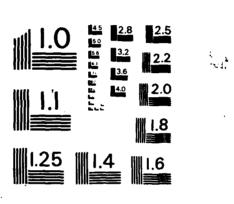
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# A ZERO EXTRACTION AND SEPARATION TECHNIQUE FOR SURFACE ACOUSTIC WAVE AND DIGITAL SIGNAL PROCESSING FIR FILTER IMPLEMENTATION

BY

KEITH V. LINDSAY B.S.E., University of Central Florida, 1984

#### THESIS

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#### **ABSTRACT**

Presented is a new method of separating the zeros of a Finite Impulse Response (FIR) filter producing an optimal digital filter or surface acoustic wave (SAW) design implementation. Overviews of zero extraction algorithms and of FIR filter design using the Remez Exchange algorithm are presented (McClellan et al. 1973).

The computer aided design (CAD) procedure presented allows the designer to specify the general filter characteristic which the Remez algorithm translates to FIR time domain coefficients. These coefficients are readily translated to the frequency (z) domain, producing an Nth order polynomial in z. The characteristic polynomial is factored to determine all roots or zeros using a three-stage factoring program presented by M.A. Jenkins (1975). The roots are optimally separated into two groups, each of which is recombined to form mutually exclusive functions. The two functions are then implemented as transducers of a SAW device or as a two-processor digital filter. The concept may be extended to more than two subgroups for multi-processor digital filter designs.

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#### **ACKNOWLEDGEMENTS**

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It is rarely known from the outset of completing a thesis what results will be achieved and the effort required to achieve those results. This effort was greatly aided by various agencies around the University of Central Florida, such as the Computer Engineering VAX facility and personnel, the Library staff, and Dr. Al Pozefsky's State Technology Applications Center, to name a few. The task of assembling this report was greatly reduced by Sharon Darling, whose professional typing skills are readily apparent.

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#### CHAPTER I

#### INTRODUCTION

Numerous techniques exist for designing and implementing finite impulse response (FIR) filters. Many of these techniques can be traced to antenna array design methods popularized in the 1940s and 1950s (Balanis 1982). Among the more prominent antenna array design techniques are the methods by Fourier transform, Schelkunoff polynomial, Dolph-Chebyshev, Taylor line-source (Chebyshev Error) and Woodward. These have spawned many of the popular contemporary FIR design techniques such as the Remez exchange algorithm based on the Chebyshev Error method, popularized by McClellan, Parks and Rabiner (1973) and non-iterative eigenfunction synthesis design, introduced by Devries (1973) and similar to Woodward's method. Another FIR design approach employs a technique known as linear programming (Rabiner 1972a,b).

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Each of these techniques will yield FIR transfer functions which may be readily implemented using a surface acoustic wave (SAW) device or a digital filter. The SAW filter, a two-transducer device, requires that the transfer function be split in some fashion between the transducers. The digital filter may be optionally implemented using two (or more) processors to increase throughput rate.

The conventional approach to implementing the SAW device is, basically, to construct one transducer such that it contains the entire FIR and to construct the other transducer such that it emulates a rect function. This imposes a requirement upon the first transducer that it be capable of handling all of the dynamics associated with the transfer function. Similarly, a single-processor digital filter implementation demands that the processor be able to handle a wide range of FIR coefficients, as well as all of them at once. These are not necessarily optimal implementations.

Some efforts to evenly split the transfer function between the two transducers of a SAW filter have been made by Morimoto, Kobayashi and Hibino (1980) and in work by Ruppel, Ehrmann-Falkenau, Stocker and Mader (1984, 1985). In both cases, these teams split the transfer function into groups of alternating zeros or roots of the transfer function about the unit circle. Any attempts to further optimize the filters were made at the expense of altering the overall frequency response in a process called compensation.

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This thesis presents an approach to near optimally split the transfer function between the two transducers or processors without altering the overall frequency response. The approach seeks to minimize non-linear and finite wordlength error effects by reducing the required tap range for each transducer, or the required range of coefficients used in a fixed-point digital processor.

#### CHAPTER II

#### OBJECTIVE OF PROPOSED WORK

# Optimal FIR Implementation Via Two-Transducer Design

Generally, the word "optimal" implies having attained a most favorable condition or degree. Many parameters must be considered during the design of a SAW or digital filter. Addressed in this thesis are those concerned primarily with filter order and coefficient dynamic range.

#### FIR Coefficients Via the Remez Algorithm

The first stage of the design is accomplished using the Remez exchange algorithm (Remez 1957) to generate a "best fit" Chebyshev polynomial to a set of frequency response specifications. The approximation, and subsequent conversion to an impulse response, is accomplished by the modified McClellan, Parks and Rabiner (1973) program presented in Appendix A. The program has been altered to permit the design of filters of up to an order of one-thousand. An initial guess of optimal filter order is obtained using a formulation presented by Vaidyanathan (1985), which states:

$$N_{e} = \frac{-10 \log_{10} \delta_{1} \delta_{2} - 13}{14.6 \Delta f}$$
 (2.1)

where:

$$\Delta f = (\omega_s - \omega_p)/2\pi$$

 $\omega_s$  = stop-band edge frequency

 $\omega_{D}$  = pass-band edge frequency

 $\delta_1$  = pass-band tolerances

 $\delta_2$  = stop-band tolerances

This  $N_{\rm e}$  provides a starting point for the filter order in the program. The program then iterates, increasing N each time, until the specifications are met.

Once the FIR coefficients are found by the modified McClellan, Parks and Rabiner (1973) program, they may be easily arranged as the coefficients of a z-domain polynomial, i.e., the z-transform of:

$$h(n) = h(t - nT) = \begin{cases} a_n & n = 0, 1, 2, ..., N \\ 0 & otherwise \end{cases}$$
 (2.2)

is

$$H(z) = \sum_{n=0}^{N} h(n)z^{-n}$$
 (2.3)

where:

 $a_n$  = the coefficients generated by the program

At this point, the impulse response could be implemented as a single SAW transducer or as a single processor digital filter. In order to split up the response between two transducers (or processors), H(z) can be expressed as:

$$H(z) = \frac{\begin{cases} \sum_{k=0}^{N} a_k z^{(N-k)} \rbrace}{z^N} = H_1(z) H_2(z)$$
 (2.4)

This equation shows all of the poles of a finite impulse response filter to be at z=0. All of the filter's zeros may be found by factoring the numerator. A judicious separation of the zeros (and poles) can then be assigned to  $H_1(z)$  and  $H_2(z)$ , the two transducers of the SAW filter. In a similar fashion, the transfer function could be split into  $H_{1...N}(z)$  for a DSP filter to increase speed and dynamic range.

Obtaining the Zeros of the Characteristic

This is the most difficult phase of the optimization. The factoring of high order polynomials was the subject of considerable effort by mathematicians during the mid-seventies. Of the many techniques surveyed, the Jenkins and Traub (1970) algorithm, and program by Jenkins (1975), appeared to be the best choice. This algorithm employs a three-stage process to determine the roots of an Nth order real polynomial. It is globally convergent and does so very rapidly. The program is extremely well written and is quite elaborate, to the point of compensating for specific machine accuracy limitations.

The Jenkins (1975) program factors the H(z) numerator polynomial and returns the real and imaginary portions of the roots of H(z). The complex roots will always appear in one of two possible ways. A set of roots may appear as complex conjugate pairs (quadratic

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factors), each with a magnitude of one corresponding to the |z|=1 unit circle. These roots will always correspond to the stopband zeros of a filter design. Another form in which they may appear is as a set of two complex conjugate pairs arranged symmetrically about the unit circle so as to satisfy the condition that:

$$z_1 z_2^* = 1$$
 (2.5)

These zeros correspond to the passband zeros of the filter. A diagram best illustrates these concepts (see Figure 1). Real roots may occur on the unit circle or in a manner similar to the passband zero case. Summarizing the above, zeros of H(z) will occur as first, second and fourth order factors.

Selection of Zeros and Subsequent Reconstitution

Once the zeros of the transfer function have been determined,
they must be separated into sub-groups and recombined. A simple
algorithm which generates all possible combinations of N roots taken
K at a time is used to separate the roots and form the sub-groups.
A polynomial is constructed from each group by multiplying the roots
within its group and the split design is rated as to the
desirability of the design.

The criteria used in the selection process employed here seeks to maximize the average tap or coefficient value, while minimizing the range and variance of those same values. These qualities are

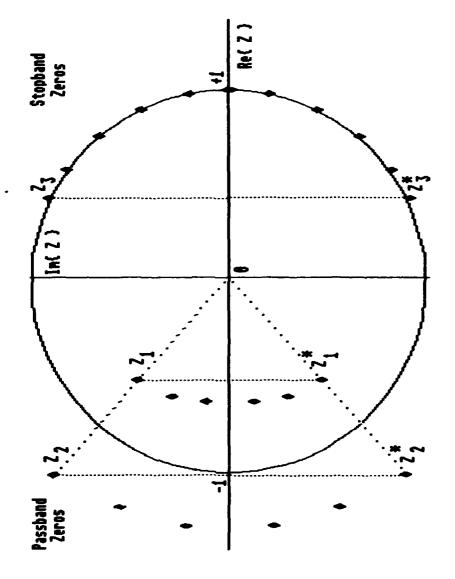


Figure 1. Typical Filter Zero Locations.

desirable in SAW devices since we usually desire a maximum finger overlap and want to avoid, as much as possible, large numbers of very small tap weights which may increase diffraction effects. In the case of digital filters, these problems translate to finite word length problems and device dynamic range.

In order to evaluate the relative merits of one combination over another, the following figure of merit is proposed:

Design FOM = 
$$\frac{\overline{x}}{R_x \sigma_x^2}$$
 (2.6)

where:

 $\overline{x}$  = average of the tap weights of both transducers combined

 $R_{\chi}$  = the range of the coefficients

 $\sigma_{x}^{2}$  = the variance of the coefficients

The ratio yields a figure of merit used to rate a given design.

Obviously, this concept could be extended to more than two

sub-transfer functions for the digital filter case.

This thesis proposes to study low order filters using this separation/reconstitution technique and to apply detected trends, if any, in a general sense.

#### CHAPTER III

#### FIR FILTER DESIGN

An excellent review of finite impulse response filter theory is presented by Lawrence R. Rabiner and Bernard Gold (1975) and by Rabiner, McClellan and Parks (1975). This review provides a theoretical background for implementing the Weighted Chebyshev Approximation filter design technique via the Remez Exchange Algorithm (Remez 1957). That presentation draws upon the work of Parks and McClellan (1973), who devised a general computer program incorporating the above. Their program was used in this thesis to provide the FIR transfer functions. An overview of the theory leading from FIR theory to a brief program description is presented. The following is adopted from Rabiner and Gold (1975).

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#### FIR Filter Frequency Response Overview

A finite impulse response describes a system which can be modeled by a difference equation in the form:

$$y(n) = \sum_{r=0}^{M} (\frac{b_r}{a_0}) x(n-r)$$
 (3.1)

Since the system output is the convolution of the system input, x(n), with the system impulse response, h(n), the impulse response can be readily seen to be:

$$h(n) = \begin{bmatrix} b_n \\ --, & n = 0, 1, 2, ..., M \\ a_0 \\ 0, & \text{otherwise} \end{bmatrix}$$
 (3.2)

The above equation describes the discrete time domain coefficients of the system FIR. This same response may be described in the frequency domain by taking the Fourier transform of h(n) to obtain  $H(e^{j\omega})$ :

$$H(e^{j\omega}) = \sum_{k=-\infty}^{\infty} h(k) e^{-j\omega k}$$
 (3.3)

Since h(n) is finite in length with respect to time,  $H(e^{j\omega})$  must be infinite with respect to frequency. However, for discrete, sampled systems,  $H(e^{j\omega})$  is periodic with respect to the sampling frequency, i.e.,:

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$$H(e^{j\omega}) = H[e^{j(\omega+2\pi m)}] \qquad m = 0, \pm 1, \pm 2, \dots$$
 (3.4)

which is periodic in frequency with a period of  $2\pi$ . This fact allows us to restrict our requirement to define  $H(e^{j\omega})$  in practical filtering applications in terms of the sampling frequency, consisting of N samples over a period equaling the length of the time domain impulse response (without any augmenting zeros). It also allows us to state that:

$$H(e^{j\omega}) = \sum_{k=0}^{N-1} h(k) e^{-j\omega k}$$
 (3.5)

The function,  $H(e^{j\omega})$ , can be described in terms of its magnitude and phase as:

$$H(e^{j\omega}) = + |H(e^{j\omega})| e^{j\theta(\omega)}$$
 (3.6)

or

$$H(e^{j\omega}) = \hat{H} (e^{j\omega}) e^{j\theta(\omega)}$$
 (3.7)

where:

 $\hat{H}$  (e<sup>j $\omega$ </sup>) = a real function

 $\theta(\omega)$  = constrained to describe a linear phase characteristic, i.e.,:

$$\theta(\omega) = -\alpha\omega \qquad -\pi \leq \omega < \pi \tag{3.8}$$

with constant phase delay implied by the constant, -  $\alpha$ . The function can be written in trigonometric form as:

$$+ |H(e^{j\omega})| \{\cos(\alpha\omega) - j \sin(\alpha\omega)\}$$
 (3.9)

In order to find  $\alpha$ , equate the real and imaginary parts. Then, we may describe  $\cos{(\alpha\omega)}$  and  $\sin{(\alpha\omega)}$  as:

$$+ |H(e^{j\omega})| \cos (\omega\omega) = \sum_{n=0}^{N-1} h(n) \cos (\omega n)$$
 (3.10)

and

$$+ |H(e^{j\omega})| \sin(\omega\omega) = \sum_{n=0}^{N-1} h(n) \sin(\omega n)$$
 (3.11)

and set up the ratio:

$$\frac{\sin (\alpha \omega)}{\cos (\alpha \omega)} = \tan (\alpha \omega) = \frac{\sum_{n=0}^{N-1} h(n) \sin (\omega n)}{N-1}$$

$$\sum_{n=0}^{N-1} h(n) \cos (\omega n)$$
(3.12)

Cross multiplying:

N-1  

$$\Sigma$$
 h(n) sin ( $\infty$ ) cos ( $n\omega$ ) -  $\Sigma$  h(n) cos ( $\infty$ ) sin ( $n\omega$ ) = 0  
n=0 (3.13)

Using the trig identity:

$$\sin (u-v) = (\sin u)(\cos v) - (\cos u)(\sin v)$$
 (3.14)

yields:

the second secon

$$\begin{array}{l} N-1 \\ \Sigma \quad h(n) \quad \sin[(\alpha-n)\omega] = 0 \\ n=0 \end{array} \tag{3.15}$$

Equation (3.15) is in the form of a Fourier series. For equation (3.15) to be valid for an odd symmetrical series (see Figure 2),  $\alpha$  and h(n) must be:

$$\alpha = \frac{N-1}{2} \tag{3.16}$$

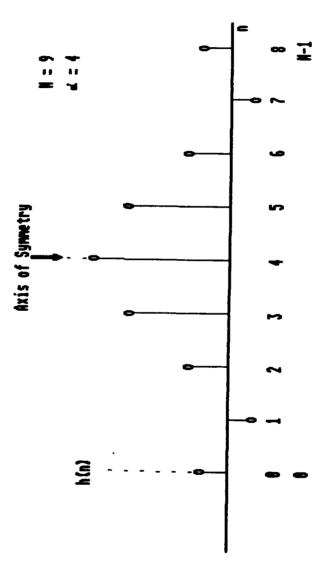


Figure 2. Odd Symmetrical FIR Series.

and:

$$h(n) = h(N-1-n)$$
  $0 \le n \le N-1$  (3.17)

In the case of an even, symmetrical series (see Figure 3),  $\alpha$  will not be an integer. The use of the fractional delay obtained here is of primary significance when designing differentiators and Hilbert transformers. These are not discussed here, but the reader is referred to Rabiner and Gold (1975) for an in-depth discussion. These values for  $\alpha$  and h(n) hold for constant group delay and constant phase filters. If constant phase delay (phase divided by frequency) is not required, i.e.,:

$$H(e^{j\omega}) = \pm |H(e^{j\omega})| e^{j\omega(\frac{\beta}{\omega} - \alpha)} = \pm |H(e^{j\omega})| e^{j(\beta - \alpha\omega)}$$
(3.18)

where the phase delay is given by  $-\frac{\beta}{\omega} + \alpha$ , then similar development (Rabiner and Gold 1975) for an odd, anti-symmetric series (see Figure 4) will lead to the result that:

$$\alpha = \frac{N-1}{2} \tag{3.19a}$$

$$\beta = \pm \frac{\pi}{2} \tag{3.19b}$$

and

$$h(n) = -h(N-1-n)$$
  $0 \le n \le N-1$  (3.19c)

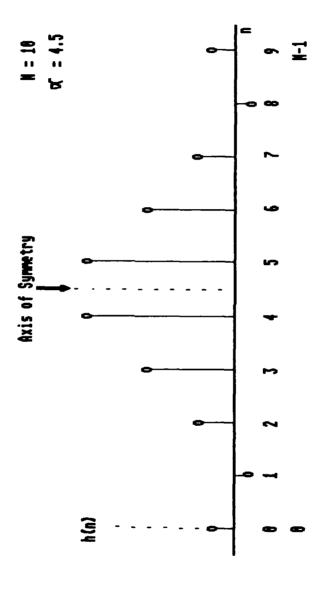


Figure 3. Even Symmetrical FIR Series.

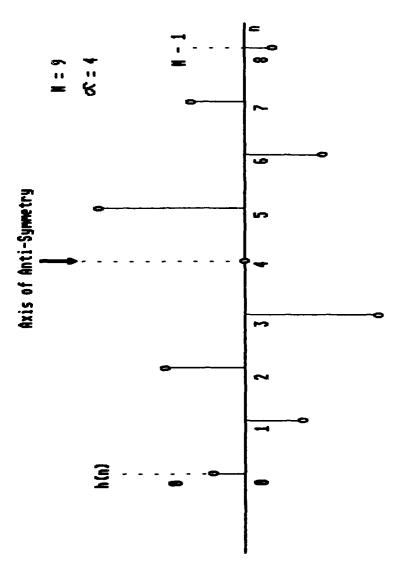


Figure 4. Odd Anti-Symmetrical FIR Series.

Again, the case of an even, anti-symmetric series (see Figure 5) is of primary interest in designing differentiators and Hilbert transformers. Equations (3.16) through (3.19) suggest four general classes that might characterize a linear phase finite impulse response filter:

- 1. Symmetrical impulse response, N odd
- 2. Symmetrical impulse response, N even
- 3. Anti-symmetrical impulse response, N odd
- 4. Anti-symmetrical impulse response, N even It is now possible to describe  $H(e^{j\omega})$  to account for these possibilities in the general relationship:

$$H(e^{j\omega}) = \hat{H}(e^{j\omega}) e^{j(\beta-\alpha\omega)}$$
 (3.20)

Rabiner and Gold (1975) next develop equations to define  $H(e^{j\omega})$  in terms of  $\hat{H}(e^{j\omega})$  for each of the above cases. The results of these developments are summarized as follows:

Case 1: Symmetrical impulse response, N odd

$$\hat{H}(e^{j\omega}) = \sum_{n=0}^{(N-1)/2} a(n) \cos(\omega n)$$
 (3.21)

with

$$a(0) = h[(N-1)/2]$$
, and 
$$a(n) = 2h[\frac{(N-1)}{2} - n] \quad \text{for } n = 1, 2, ..., (N-1)/2$$

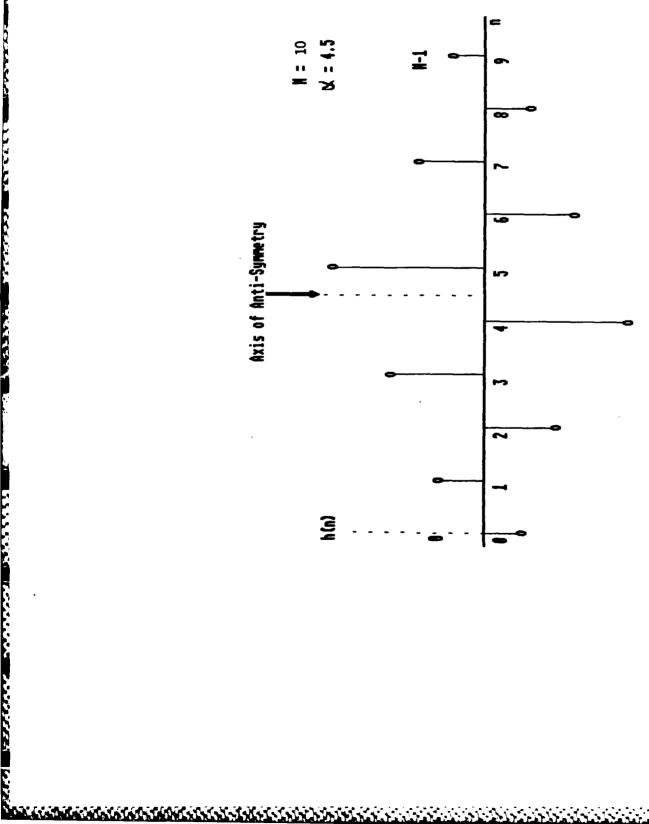


Figure 5. Even Anti-Symmetrical FIR Series.

which yields:

$$H(e^{j\omega}) = e^{-j\omega(N-1)/2}$$
  $\sum_{n=0}^{(N-1)/2} a(n) \cos(\omega n)$  (3.23)

Case 2: Symmetrical impulse response, N even

$$\hat{H}(e^{j\omega}) = \sum_{n=1}^{N/2} b(n) \cos \left[\omega(n-0.5)\right]$$
 (3.24)

with

$$b(n) = 2h(N/2 - n), \quad n = 1, 2, ..., N/2$$
 (3.25)

$$H(e^{j\omega}) = e^{-j\omega(N-1)/2} \sum_{n=1}^{N/2} b(n) \cos [\omega(n-0.5)]$$
 (3.26)

Case 3: Anti-symmetrical impulse response, N odd

$$\hat{H}(e^{j\omega}) = \sum_{n=1}^{(N-1)/2} c(n) \sin(\omega n)$$
 (3.27)

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$$c(n) = 2h[(N-1)/2 - n], \quad n = 1, 2, ..., (N-1)/2$$
 (3.28)

$$H(e^{j\omega}) = e^{-j\omega(N-1)/2} e^{j\pi/2} \int_{n=1}^{(N-1)/2} c(n) \sin(\omega n)$$
 (3.29)

Case 4: Anti-symmetrical impulse response, N even

$$\hat{H}(e^{j\omega}) = \sum_{n=1}^{N/2} d(n) \sin \left[\omega(n-0.5)\right]$$
 (3.30)

with

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$$d(n) = 2h(N/2 - n), \qquad n = 1, 2, ..., N/2$$
 (3.31)

$$H(e^{j\omega}) = e^{-j\omega(N-1)/2} e^{j\pi/2} \int_{n=1}^{N/2} d(n) \sin [\omega(n-0.5)] (3.32)$$

#### Weighted Chebyshev Approximation

The frequency response of the desired system is completely described by  $H(e^{j\omega})$ . From the development in the first section of this chapter, it is obvious that this description for each case can be considered as a series of sine or cosine functions. These series can be easily related to Chebyshev polynomials.

The Chebyshev polynomial represents an expansion of cos (mu) for any value of m. We know that any real function can be represented as a sum of sinusoids. These sinusoids are of the form cos (mu), with m indicating the highest harmonic required to reconstruct the original function. Of course, some functions require that m approach  $\infty$ . Within given limits, however, it is possible to represent a desired frequency response curve as a sum of Chebyshev polynomials of finite length m. The Chebyshev polynomial expansions take the following forms:

$$\frac{m}{0}$$
  $\frac{\cos mu}{1}$  = 1  
1  $\cos u = \cos u$   
2  $\cos 2u = 2\cos^2 u - 1$   
3  $\cos 3u = 4\cos^3 u - 3\cos u$ 

Letting  $z = \cos(u)$ , or  $u = \cos^{-1}(z)$ , then:

<u>m</u>	cos mu	Chebyshev <u>Designation</u>	(3.34)
0	1	T <sub>0</sub> (z)	
1	Z	T <sub>1</sub> (z)	
2	$2z^2 - 1$	T <sub>2</sub> (z)	
3	4z <sup>3</sup> - 3z	T <sub>3</sub> (z)	

A recursive relationship emerges:

$$T_{m}(z) = \cos [m \cos^{-1} (z)] = \cos (mu), -1 \le z \le 1$$
 (3.35)

In essence, we are using a sum of Chebyshev polynomials to curve-fit to the desired frequency response from zero to one-half the sampling frequency. Given the four cases discussed at the end of the first section of this chapter, and using Chebyshev polynomials to represent the series, the problem defaults to determining the

scaling of the coefficients with respect to frequency. This process is called "weighting" the approximation and is discussed in depth by Rabiner and Gold (1975). It is reiterated here briefly.

For the four cases described in the first section of this chapter, a general expression can be written to define  $H(e^{j\omega})$  as:

$$H(e^{j\omega}) = e^{-j\omega(N-1)/2} e^{j(\pi/2)L} \hat{H}(e^{j\omega})$$
 (3.36)

The exponent L will take on a value of either 0 or 1, depending upon the case considered. Now, a table can be constructed which shows values for L and the form of  $\hat{H}(e^{j\omega})$  for the appropriate case of symmetry and N (see Table 1). The previous discussion of the form of the Chebyshev polynomial would suggest that the expressions for  $\hat{H}(e^{j\omega})$  may be converted to summations involving cosines (as opposed to sines) using ordinary trigonometric identities. Once this is done, Table 1 can be rewritten in terms of functions which are fixed functions of  $\omega$ , which will be referred to as  $Q(e^{j\omega})$ , and as functions of the cosine series, which will be referred to as  $P(e^{j\omega})$  (see Table 2). For cases 2 through 4,  $Q(e^{j\omega})$  is constrained to be zero at either  $\omega = 0$  or  $\omega = \pi$ , or both.

Now, it is possible to set up a relationship between the desired response at given frequencies to within a prescribed accuracy. To do so, let  $D(e^{j\omega})$  represent the desired response of the filter and let  $W(e^{j\omega})$  represent the weighting on the allowable error as a function of frequency regions or bands (i.e., the ratio of the

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TABLE 1
DEFINITION OF L FOR THE FOUR FILTER CASES

	٦	н(е <sup>јω</sup> )
Case 1: Symmetrical impulse response, N odd	0	$\hat{H}(e^{j\omega}) = \frac{(N-1)/2}{\Sigma}$ a(n) cos (ωn) n=0
Case 2: Symmetrical impulse response, N even	0	$\hat{H}(e^{j\omega}) = \sum_{n=1}^{N/2} b(n) \cos [\omega(n-0.5)]$
Case 3: Anti-symmetrical impulse response, N odd	-	$\hat{H}(e^{j\omega}) = \sum_{n=1}^{\infty} c(n) \sin(\omega n)$
Case 4: Anti-symmetrical impulse response, N even	1	$\hat{H}(e^{j\omega}) = \sum_{n=1}^{N/2} d(n) \sin [\omega(n-0.5)]$

TABLE 2 Q(e^j  $^\omega$ ) AND P(e^j  $^\omega$ ) DEFINED FOR THE FOUR FILTER CASES

govern commence exercises, normally exercises, sometimes and exercises. Sometimes

	0(e <sup>Ĵω</sup> )	P(e <sup>jω</sup> )
Case 1: Symmetrical impulse response, N odd	<b></b> 1	$\hat{H}(e^{j\omega}) = \sum_{n=0}^{(N-1)/2} a(n) \cos(\omega n)$
Case 2: Symmetrical impulse response, N even	( <mark>∂</mark> ) soo	$\hat{H}(e^{j\omega}) = \sum_{n=0}^{\infty} b(n) \cos(\omega n)$
Case 3: Anti-symmetrical impulse response, N odd	sin (ω)	$\hat{H}(e^{j\omega}) = \sum_{n=0}^{\infty} c(n) \cos(\omega n)$
Case 4: Anti-symmetrical impulse response, N even	sin (½)	$\hat{H}(e^{j\omega}) = \sum_{n=0}^{\Sigma} d(n) \cos(\omega n)$

stopband ripple to the passband ripple). With these functions, the error for a given approximation can be calculated as:

$$E(e^{j\omega}) = W(e^{j\omega}) [D(e^{j\omega}) - \hat{H}(e^{j\omega})]$$
 (3.37)

with  $\hat{H}(e^{j\omega})$  being the trial design.  $\hat{H}(e^{j\omega})$  can be separated into its two symbolic parts to yield:

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$$E(e^{j\omega}) = W(e^{j\omega}) [D(e^{j\omega}) - P(e^{j\omega}) Q(e^{j\omega})]$$
 (3.38)

 $Q(e^{j\omega})$  may be factored out of the quantity in parentheses since it is a fixed function of frequency. This yields:

$$E(e^{j\omega}) = W(e^{j\omega}) Q(e^{j\omega}) [D(e^{j\omega})/Q(e^{j\omega}) - P(e^{j\omega})] \qquad (3.39)$$

Defining  $[W(e^{j\omega}) Q(e^{j\omega})]$  as  $\widehat{W}(e^{j\omega})$ , and  $[D(e^{j\omega})/Q(e^{j\omega})]$  as  $\widehat{D}(e^{j\omega})$ , equation (3.39) may be rewritten as:

$$E(e^{j\omega}) = \hat{W}(e^{j\omega}) [\hat{D}(e^{j\omega}) - P(e^{j\omega})]$$
 (3.40)

The problem now defaults to finding the values for the coefficients of the Chebyshev polynomials  $[P(e^{j\omega})]$  such that the maximum error over each specified frequency band is minimized. To accomplish this, the Alternation Theorem is used. It states (Rabiner and Gold 1975):

Theorem: If  $P(e^{j\omega})$  is a linear combination of r cosine functions, i.e.,:

$$P(e^{j\omega}) = \sum_{n=0}^{r-1} \alpha(n) \cos(n\omega)$$
 (3.41)

then a necessary and sufficient condition that  $P(e^{j\omega})$  be the unique, best weighted Chebyshev approximation to a continuous function  $\widehat{D}(e^{j\omega})$  on A, a compact subset of  $(0,\pi)$ , is that the weighted error function  $E(e^{j\omega})$  exhibit at least (r+1) extremal frequencies in A; i.e., there must exist (r+1) points  $\omega_i$  in A such that  $\omega_1 < \omega_2 < \ldots < \omega_{r+1}$  and such that  $E(e^{j\omega i}) = -E(e^{j\omega i+1})$ ,  $i = 1, 2, \ldots, r$ , and  $|E(e^{j\omega i})| = \max [E(e^{j\omega})]$  for all  $\omega$  in A.

Rabiner and Gold (1975) show that for the four cases of filter design presented, that the number of extremal frequencies in  $\hat{H}(e^{j\omega})$  obey the following constraints:

Case 1: 
$$N_e \le (N+1)/2$$
  
Case 2:  $N_e \le N/2$   
Case 3:  $N_e \le (N-1)/2$   
Case 4:  $N_e \le N/2$ 

The extremal frequencies, or extrema, are divided up between the stop and passbands of the filter and they describe the peaks and troughs of the Chebyshev approximation. A diagram best describes the relationship of the extremal frequencies and the shape of the Chebyshev approximation waveform (see Figure 6).

There are several ways in which to obtain the extremal frequencies of  $\hat{H}(e^{j\omega})$ . The first method, described briefly, was originally proposed by Herrmann and Schuessler. The method capitalizes on the fact that a local maxima (+8) or a local minima

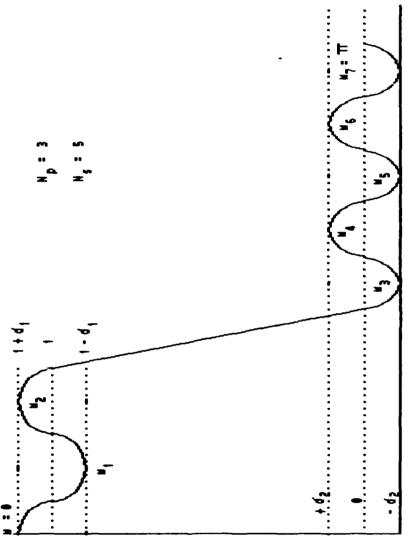


Figure 6. Chebyshev Approximation Extremal Frequencies.

 $(-\delta)$  occurs in the region of an extrema, and that the derivative is zero at that point. Two equations in N<sub>e</sub> with two N<sub>e</sub> unknowns  $(N_e$  impulse response coefficients and N<sub>e</sub> frequencies where  $\hat{H}(e^{j\omega})$  obtains an extremal value) are:

$$\hat{H}(e^{j\omega_i}) = \pm \frac{\delta}{j\omega_i} + D(e^{j\omega_i})$$
  $i = 1, 2, ..., N_e$  (3.43)

and

$$\frac{d \left(\hat{H}(e^{j\omega})\right)}{d\omega} = 0 \qquad i = 1, 2, ..., N_e \qquad (3.44)$$

where  $E(e^{j\omega i}) = \pm \delta$ , and these are solved iteratively for values of  $N_e$ . This procedure works well for filters with an order about 60 or less.

Another method used is one devised by Hofstetter, Oppenheim and Siegel which is called the Polynomial Interpolation Solution by Rabiner and Gold (1975). The basic idea behind this algorithm is that an initial guess of the extremal frequencies is made and  $\hat{H}(e^{j\omega})$  is evaluated at these points. The algorithm then searches for the actual extrema found during that trial and iterates again, this time using the newly found extrema. Eventually, the process converges to the minimum ripple attainable for a given  $N_e$ . Very large order filters can be designed by this method. The Polynomial Interpolation Solution technique is very similar to the last technique to be described, the Remez Exchange Algorithm.

# The Remez Exchange Algorithm

We have seen that the goal of obtaining the desired response  $\hat{D}(e^{j\omega})$  is met by obtaining the approximating function  $P(e^{j\omega})$  which best minimizes the weighted error function  $E(e^{j\omega})$ . The Remez Exchange Algorithm accomplishes this by using a dense grid of frequency points to find the extremal frequencies. An initial guess as to the location of the (r+1) frequencies is made, similar to the Polynomial Interpolation Solution method. Then, the error function is forced to have a value of  $\underline{+}\ \delta$ . The signs alternate, since the extrema are expected to alternate above and below the indicated level by  $\delta$  in the final design. These constraints generate the separate error function  $[E(e^{j\omega})]$  equation for each extremal frequency, given from equation (3.40) as:

$$\hat{W}(e^{j\omega_k})[\hat{D}(e^{j\omega_k}) - P(e^{j\omega_k})] = (-1)^k \delta$$
  $k = 0, 1, ..., r$  (3.45)

which generates an (r+1) x (r+1) matrix of equations to solve. Remez (1957) found an alternative closed-form solution (appropriately modified for the current variable set) to be:

$$\delta = \frac{a_0 \hat{D}(e^{j\omega_0}) + a_1 \hat{D}(e^{j\omega_1}) + \dots + a_r \hat{D}(e^{j\omega_r})}{a_0 / \hat{W}(e^{j\omega_0}) - a_1 / \hat{W}(e^{j\omega_1}) + \dots + (-1)^r a_r / \hat{W}(e^{j\omega_r})}$$
(3.46)

where:

$$a_k = \prod_{\substack{i=0\\i\neq k}}^{r} \frac{1}{(x_k - x_i)}$$
 (3.47)

and

$$x_{j} = \cos \omega_{j} \tag{3.48}$$

At this point, the optimum  $\delta$  for a given set of extremal frequencies is known. The next step is to form the approximating function  $P(e^{j\omega})$  along the r extrema points by using the barycentric form of the Lagrange interpolation formula:

$$P(e^{j\omega}) = \frac{\sum_{k=0}^{r-1} \left(\frac{\beta_k}{x - x_k}\right) c_k}{\sum_{k=0}^{r-1} \frac{\beta_k}{(x - x_k)}}$$
(3.49)

where:

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$$\beta_{k} = \prod_{\substack{i=0\\i\neq k}}^{r-1} \frac{1}{(x_{k} - x_{i})}$$
 (3.50)

and

$$C_k = \hat{D}(e^{j\omega_k}) - (-1)^k \frac{\delta}{\hat{\omega}(e^{j\omega_k})} \quad k = 0, 1, ..., r-1$$
 (3.51)

$$x_i = \cos \omega_i$$

$$x_k = \cos \omega_k$$

$$x = \cos \omega$$
(3.52)

Once the approximating function  $P(e^{j\omega})$  has been formed, it is possible to evaluate  $E(e^{j\omega})$  along a dense set of frequencies which are equally spaced along the frequency axis from zero to one-half the sample frequency. If:

$$|\mathsf{E}(\mathsf{e}^{\mathsf{j}\omega})| \leq \delta$$
 (3.53)

then an optimal approximation to the desired frequency response has been found. If the weighted error function exceeds  $\delta$ , then a new set of (r+1) extremal frequencies is chosen by selecting the peaks of the error curve. This process quickly forces  $\delta$  to converge to its maximum value for a given number of extremal frequencies. If there are more than (r+1) extrema in  $E(e^{j\omega})$ , then the new number of extrema is retained and used in the next iteration of the process.

The final impulse response coefficients are obtained by performing a  $2^M$  point Inverse Discrete Fourier Transform on  $P(e^{j\omega})$ , where  $2^M \ge N$ . Note that this N is the filter order plus 1.

#### CHAPTER IV

## ZERO EXTRACTION TECHNIQUES

The problem of extracting the zeros of a polynomial turns out to be far from simple, sparking the interest of mathematicians and scientists for centuries. With the advent of the digital computer, the factoring of high order polynomials has become possible, though not entirely without grief. Some of the more prominent approaches and associated problems are briefly discussed here.

# Polynomial Theory

A polynomial in z is an equation which takes the form:

$$a_N z^N + a_{N-1} z^{N-1} + \dots + a_2 z^2 + a_1 z + a_0$$
 (4.1)

or, alternatively,

$$\sum_{n=0}^{N} a_n z^n \tag{4.2}$$

where the coefficients  $a_N$ ,  $a_{N-1}$ , ...,  $a_0$  are real numbered constants. This form of equation is readily identified with the equation describing the finite impulse response filter z-domain representation. This polynomial can be expressed also as a product of its roots or zeros as:

$$\prod_{n=1}^{N} (z - z_n)$$
(4.3)

Notice that for a polynomial with N roots, there are N product-form terms and N+1 summation-form terms. This can cause some confusion at times. For example, the McClellan, Parks and Rabiner (1973) program discussed in the previous chapter displays a filter order of N when, in fact, N coefficients are actually meant.

# Zero Characteristics of FIR Filters

When described by a polynomial in the  $z=e^{j\omega}$  plane, FIR filter zeros plotted in the z-plane always have a distinct appearance. All of the stopband zeros will occur exactly on the unit circle and will always occur as complex conjugate pairs, unless they are real. The complex passband zeros will always occur in sets of four, one inside the unit circle, another outside such that the magnitude of one multiplied by the other will equal exactly one. This pair also has corresponding complex conjugates, hence the set of four.

Due to the nature of the Chebyshev polynomial approximation of the FIR frequency response, there are no repeating zeros or multiple roots to contend with. However, this does pose problems in other factoring situations and is discussed briefly.

# Factoring Methods

Several unique approaches to zero extraction exist. The first was by none other than Sir Issac Newton (1642-1727). Since that

time, other algorithms by Bairstow, Lin, Muller and Birge-Vieta have arrived (Ralston and Wilf 1960). These algorithms have the relative liability of not being able to assure convergence for any initial guess of a root. Other methods which virtually assure convergence within a class of problems are methods of Lehmer, Graeffe and Bernoulli. The Bisection method is probably the most crude method of root-finding, relying upon a purely iterative process of testing discrete points in the z-plane until the roots are found. The latter four cases have the disadvantage of slow convergence. There exist several matrix-oriented computer program packages, such as EISPACK (Smith et al. 1976), which are designed to find the eigenvalues of a matrix, another means of finding the zeros. However, these programs are extremely memory-inefficient. The method adopted for this thesis project is the Jenkins and Traub (1970) method which combines many of the above techniques, as well as ones by Traub, into a very complex algorithm and computer program which is convergent for a wide class of problems and is very machine-efficient. Discussed briefly are the more prominent methods.

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#### Newton's Method

The process of finding a root by Newton's method is perhaps the best known and most easily understood (Blomquist 1968). It is based on beginning with an initial guess for the root and iteratively converging to the actual root with the method of steepest descent. The following relationship describes the method:

$$z_{n+1} = z_n - P(z_n)/P^1(z_n)$$
 (4.4)

where  $z_n$  is the current guess of the root,  $P(z_n)$  is the value of the polynomial at  $z_n$ ,  $P^1(z_n)$  is the value of the derivative of  $P(z_n)$  and  $z_{n+1}$  is the next guess (or, eventually, the root). This process may be carried out iteratively to any practical, desired accuracy. The root is obtained when the difference between  $z_{n+1}$  and  $z_n$  is less than the required accuracy. Figure 7 graphically shows how successive iterations ultimately converge to the root.

Problems with this technique occur since it has no direct way of determining if a multiple root exists and the method does not always converge to the root. An example of how the method may fail is shown in Figure 8. This case demonstrates that choosing an initial guess too far from the actual root may prevent convergence. In this example, the method will oscillate between  $z_1$  and  $z_0$ , since each point represents the other's successive approximation. The method also requires the use of complex arithmetic in evaluating  $P(z_n)$  and  $P^1(z_n)$  in the case of complex roots, which further limits this method.

#### Bairstow's Method

Prior to the Jenkins and Traub (1970) approach, the Bairstow method was regarded as one of the best techniques for extracting zeros. The primary advantage which this method has over Newton's Method is that it uses only real arithmetic to evaluate the

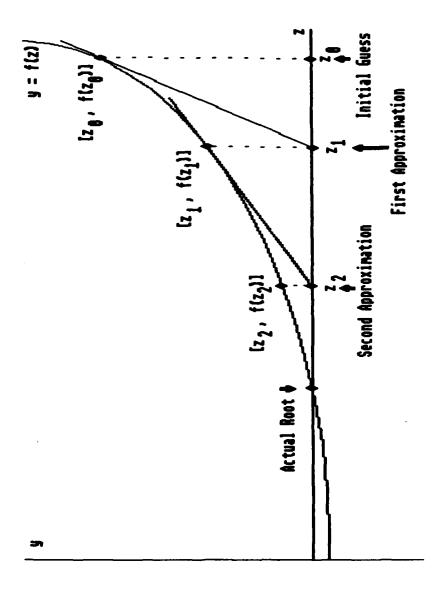


Figure 7. Newton's Method of Successive Approximation.

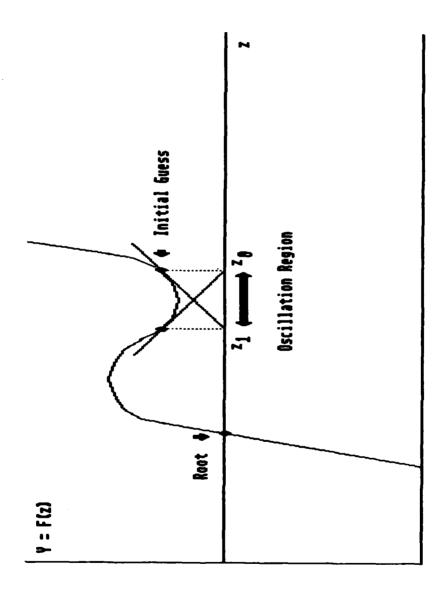


Figure 8. Newton's Method - Failure to Converge.

polynomial. The basic idea is to use Newton's Method to find unique quadratic factors to the polynomial using only real arithmetic, remove the quadratic via synthetic division and use the well-known quadratic formula to extract the complex roots of the quadratic factor. This technique automatically locates multiple roots, since it deflates the polynomial by an order of two for each quadratic factor found.

Simons, Weeks and Kotick (1983) developed an elegant formulation which best expresses Bairstow's Method. The algorithm begins with a polynomial in the form of:

$$P_N(z_n) = a_N z^N + a_{N-1} z^{N-1} + a_{N-2} z^{N-2} + \dots + a_1 z^1 + a_0 z^0$$
 (4.5)

Newton's Method is used to approach the root by using equation (4.4). However, Bairstow's Method evaluates  $P_N(z)$  and its derivative(s) at the pair of complex points  $z_n$  and  $z_n^*$  by using only real arithmetic as follows:

Let 
$$P_N(z) = \sum_{n=0}^{N} a_n z^n$$
 (4.6)

Let the quadratic factor take the form:

$$z^2 + \alpha z + \beta \tag{4.7}$$

where:

and

$$\beta = \sigma^2 + \omega^2$$

Then, dividing  $P_N(z)$  by this quadratic factor yields a polynomial of order N-2 with a remainder  $R_1z+R_0$ , i.e.,

$$\frac{P_N(z)}{z^2 + \alpha z + \beta} = P_{N-2}(z) + \frac{R_1 z + R_0}{z^2 + \alpha z + \beta}$$
 (4.8)

Multiplying both sides by the quadratic factor yields:

$$P_N(z) = P_{N-2}(z) (z^2 + \alpha z + \beta) + R_1^z + R_0$$
 (4.9)

Obviously, if  $R_1z + R_0 = 0$ , then the roots are:

$$\frac{-\alpha + \sqrt{\alpha^2 - 4\beta}}{2} \tag{4.10}$$

by the quadratic formula. Therefore, the problem defaults to choosing values for  $\alpha$  and  $\beta$  such that the remainder is zero.  $R_1$  and  $R_0$  can be related to  $P_N(z)$  by the following:

$$z^{2} + \alpha z + \beta \overline{)a_{N}z^{N-2} + (a_{N-1}-\alpha a_{N-1})z^{N-3} + [a_{N-2}-\alpha a_{N-1}+(\alpha^{2}+\beta)a_{N}]z^{N-3} + \dots}}$$

$$\underline{a_{N}z^{N} + a_{N-1}z^{N-1} + a_{N-2}z^{N-2} + a_{N-3}z^{N-3} + \dots}}$$

$$\underline{a_{N}z^{N} + \alpha a_{N}z^{N-1} + \beta a_{N}z^{N-2}}$$

$$\underline{(a_{N-1}-\alpha a_{N})z^{N-1} + (a_{N-2}-\beta a_{N})z^{N-2} + a_{N-3}z^{N-3} + \dots}}$$

$$\underline{(a_{N-1}-\alpha a_{N})z^{N-1} + (\alpha a_{N-1}-\alpha a_{N})z^{N-2} + (\beta a_{N-1}-\alpha \beta a_{N})z^{N-3}}}$$

$$\underline{Ia_{N-2}-\alpha a_{N-1} + (\alpha^{2}-\beta)a_{N}Jz^{N-2}}$$

$$+ (a_{N-3}-\beta a_{N-1}+\alpha \beta a_{N})z^{N-3} + \dots$$

$$\vdots$$

$$(4.11)$$

Letting:

$$b_{N-2} = a_{N}$$

$$b_{N-3} = a_{N-1} - a_{N}\alpha = a_{N-1} - b_{N-2}$$

$$b_{N-4} = a_{N-2} - a_{N}\alpha - a_{N-1}\alpha - a_{N}\alpha$$

$$= a_{N-2} - b_{N-2} - b_{N-3}\alpha$$
(4.12)

a recursive relationship emerges:

$$b_{N-2-n} = a_{N-2-n} - \beta b_{N-2-n} - \alpha b_{N-3-n}$$
 (4.13)

Iterating to n = N yields:

$$b_{0} = a_{2} - \beta b_{2} - \alpha b_{1}$$

$$b_{-1} = a_{1} - \beta b_{1} - \alpha b_{0}$$

$$b_{-2} = a_{0} - \beta b_{0} - \alpha b_{-1}$$
(4.14)

Since

$$\frac{R_1 z + R_0}{z^2 + \alpha z + \beta} = b_{-1} z^{-1} + b_{-2} z^{-2}$$
 (4.15)

then

$$R_{1}z + R_{0} = (z^{2} + \alpha z + \beta) (b_{-1}z^{-1} + b_{-2}z^{-2})$$
 (4.16)

$$= b_{-1}z + b_{-2} + \alpha b_{-1} + \frac{\alpha b_{-2}}{z} + \frac{\beta b_{-1}}{z} + \frac{\beta b_{-2}}{z^2}$$
 (4.17)

Equating like coefficients:

$$R_{1} = b_{-1}$$

$$R_{0} = b_{-2} + \alpha b_{-1}$$

$$= a_{0} - \beta b_{0} - \alpha b_{-1} + \alpha b_{-1}$$

$$= a_{0} - \beta b_{0}$$
(4.18)

The polynomial  $P_N(z)$  and its derivative(s) may now be evaluated at any pair of complex conjugate points specified by the chosen quadratic

factor. Newton's Method now follows easily, unless  $P_N^{-1}(z)$  evaluates to zero. If this is the case, multiple roots exist for the chosen pair,  $\sigma + j\omega$ . In this case:

$$\frac{P_{N}(z)}{P_{N}^{1}(z)} \tag{4.19}$$

is replaced by successive derivatives

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$$\frac{P_N^{M}(z)}{P_N^{M+1}(z)} \tag{4.20}$$

until  $P_N^{M+1}(z_n) \neq 0$ . Now the root, as well as its multiplicity, is known.

The final step of the process is to remove each quadratic factor from the polynomial (polynomial deflation) using synthetic division. Bairstow's Method is again applied to the resulting polynomial until all of the roots are found.

The disadvantage of Bairstow's Method lies in the necessity to make a reasonably accurate guess of the quadratic factor. A bad guess can prevent convergence in the same manner as happens in Newton's Method. Both of the above processes suffer from machine rounding problems which occur with successive deflation of the original, high ordered polynomial.

### Jenkins and Traub Method

This is a highly complex method which attempts to incorporate all of the above advantages while avoiding the mentioned pitfalls. The program incorporating the algorithm was the subject of Jenkins' doctoral dissertation (Jenkins 1975). The process incorporates three stages of zero extraction using Bairstow's Method, Newton's Method and three shifting techniques used to hasten convergence. The method assures rapid convergence for a wide class of polynomials. Zeros are removed in roughly increasing order of modulus; i.e., the zeros closest to the origin are generally removed first, the ones furthest from the origin are removed last. This is done in order to reduce the instability problems which may accompany the deflation process. A discussion of the variable shift algorithm is beyond the scope of this thesis and the reader is referred to the Jenkins and Traub (1970) paper for a formal theoretical treatment.

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One of the interesting aspects addressed by Jenkins in writing the program is that it takes into account the specific capabilities and limitations of floating point manipulations on a given machine. This feature allows the program to be customized to a specific machine in order to achieve the highest possible zero extraction accuracy for that machine (using the Jenkins and Traub algorithm).

The program in Appendix B appears to be the current state-ofthe-art in non-matrix methods of polynomial factoring. For example, the 1986 IMSL math libraries make use of this algorithm for their zero extraction approach. Schelin (1983) also indicated that this was the prime non-matrix type algorithm as of 1982.

In practice, the program handles up to roughly an order of one-hundred with little difficulty. Convergence problems begin to occur with increasing frequency beyond this limit, based upon actual tests.

# CHAPTER IV

### ZERO SEPARATION AND RECONSTITUTION

The prime reason for judicious zero separation is based on a desire to increase the dynamic range of the transducer or DSP device without sacrificing any of its transfer characteristics. The dynamic range is largely affected by relatively small FIR coefficients. These small coefficients correspond to small area overlaps of fingers in SAW devices and to small register coefficients in DSP filters. The net effect in the SAW device is for the small overlap to appear more like a point source wave generator as opposed to a desired planar wave source. Second order effects also begin to become more predominant for this situation. Similarly, DSP filters suffer from rounding effects when forced to sum products of very large and very small filter coefficients, even when floating-point arithmetic is used. The following DSP example demonstrates this problem.

- Let the internal number representation be from 1.00 to 1.00.
- 2. Let the system function be:

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$$y(n) = (1.00) x(n) + (0.02) x(n-1)$$

which is an FIR filter with coefficients 1.00 and 0.02.

3. Let x(0) = x(1) = 0.20 and x(-1) = 0.0, then it follows that:

$$y(0) = (1.00) (0.20) + (0.02) (0.00) = 0.20$$

$$y(1) = (1.00) (0.20) + (0.02) (0.20) = 0.204 (actual)$$

However, y(1) = 0.204 will be truncated to 0.20, since the internal precision only allows two decimal places.

In practice, it may not be possible to completely avoid the problems associated with dynamic range, but it is desirable to attain the best dynamic range for a given design. The above example demonstrates that three qualities of the FIR coefficients bear close scrutiny during the design process: the average, the variance and the range of the coefficient values.

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# Statistical Qualities

The highest average value for the FIR coefficients provides the greatest average finger overlap on a SAW device, thus the highest average energy injection into (or removal from) the substrate. Similarly, the highest average coefficient value in a DSP filter produces data with the highest average value within the register working ranges. Since sign changes may be handled easily in either type of device, the average is taken of the absolute value of the FIR coefficients. The average is defined as:

$$\frac{\Delta}{x} \stackrel{\Sigma}{=} \frac{|h(n)|}{N}$$
 (5.1)

The variance of the FIR coefficients indicates how much they change on the average. In other words, the variance indicates how smooth the overall envelope is. Again, since sign changes may be handled easily in either type of device, the variance is taken of the absolute value of the FIR coefficients. The variance is defined as:

$$\sigma^{2} = \frac{\sum_{n=1}^{N} [h(n)]^{2} - \{\sum_{n=1}^{N} |h(n)|\}^{2}}{N(N-1)}$$
(5.2)

The range of the coefficients is an absolute indication of now far apart the minimum and maximum coefficient values are. Since the coefficients are normally scaled to a maximum of 1.0, the range will provide an indication of the minimum coefficient value. As in the above two statistical qualities, the range is taken for the absolute coefficient values. It is defined as:

Range = 
$$|h(n)_{max}| - |h(n)_{min}|$$
 (5.3)

These three qualities allow the designer some means of rating one given design against another quantitatively, leading to a design Figure of Merit (FOM).

# The Figure of Merit (FOM)

A design objective which requires the best split of the zeros of the transfer function requires that the designer be able to rate

one design over another until the best is found. This can be done by assigning an FOM to the design based on the three statistical qualities discussed above. The goal addressed here is to split the zeros between two transducers or processors such that a gain in available dynamic range is found. To do this, the following procedure relating the combined transducer coefficient average, variance and range values to a figure of merit is proposed:

- 1. Normalize all  $h_1(n)$  coefficients to 1.0 maximum.
- 2. Normalize all  $h_2(n)$  coefficients to 1.0 maximum.

(NOTE: For the special case where all of the FIR coefficients are implemented by h<sub>1</sub>(n), then h<sub>2</sub>(n) is set to 1 to represent the impulse function since the convolution of the impulse response with an impulse is the impulse response.)

- 3. Form an array consisting of  $|h_1(n)|$ .
- 4. Concatenate the  $|h_2(n)|$  values to this array.
- 5. Find the average  $(\overline{x})$ , variance  $(\sigma^2)$  and range of the newly-formed array.
- 6. Apply the relation:

$$FOM = \frac{\overline{x}}{\sigma^2 \cdot Range}$$
 (5.4)

This equation reflects a desire to maximize the average while minimizing the variance and range of the coefficient values. It provides the designer with a means of rating one set of split zeros versus another.

## Splitting the Zeros

In order to determine if an optimum zero-splitting pattern might exist, all possible combinations for distributing the zeros must be made and each combination tested using the FOM procedure. This process is very tedious but may be readily implemented on a computer due to the iterative nature of the process. Since there are:

$$_{K=0}^{N/2} \frac{N!}{K! (N-K)!}$$
 (5.5)

total, non-repeating combinations, the process has a practical computational upper limit of about FIR order N = 40 (which has over  $10^{11}$  combinations) on a VAX 11-750. However, equations of low order may be used to model any trends applicable to the higher order cases encountered in SAW devices.

# Computer Implementation

Appendix C shows a listing of the program COMBO used to generate (Beckenbach 1964) and rate all of the combinations of zeros provided by the Jenkins (1975) program. These zeros are the result of factoring the frequency response provided by the McClellan, Parks and Rabiner (1973) program.

COMBO first generates an array consisting of all of the zeros (complex conjugate pairs or reals) to be placed in  $H_1(z)$ , while all others are assumed to be placed in  $H_2(z)$ . The program insures that

for responses which contain five real zeros, a sufficient number of combinations occur. Once a combination has been specified, program control is passed to the zero reconstitution subroutine.

The reconstitution subroutine POLYRECON is responsible for multiplying the appropriate zeros together to form a test case  $H_1(z)$  and  $H_2(z)$ . This algorithm uses either a real root to form a linear factor or a complex conjugate pair to form a quadratic factor. All arithmetic performed is real. The  $H_1(z)$  and  $H_2(z)$  arrays are readily converted to scaled  $h_1(n)$  and  $h_2(n)$  arrays. The arrays are combined as described in the FOM procedure, the statistics are performed by the HSTAT subroutine and an FOM is assigned. Next, the FOM is compared to the FOM of the previous design and a decision is made as to which is best. If the new design is best, those results are stored and the program proceeds to iterate again. Upon completion, the results are passed back to the calling routine for further analysis.

### CHAPTER VI

### COMPUTER AIDED DESIGN APPLICATION

The Solid State Devices Lab group at the University of Central Florida, headed by Dr. Donald Malocha, uses a computer analysis system called SAWCAD, developed at UCF, to design SAW filter devices. Added to this program are the McClellan, Parks and Rabiner (1973) (Remez) program, the Jenkins (1975) zero location program and the Optimizing program discussed in Chapter V. The following is a brief discussion of the use of the added features to SAWCAD. The complete SAWCAD package is not discussed here. Further information on other aspects of SAWCAD can be obtained by referring to Richie (1983).

A listing of the main menu is shown in Figure 9. A filter design is initiated by selecting the (C)hebyshev [REMEZ] option to the main menu. The program proceeds to the McClellan, Parks and Rabiner (1973) program which has been somewhat modified. The data entry routine consists of an interactive session between the computer and the designer. During this session, the designer is asked to specify the filter type (i.e., bandpass, differentiator or Hilbert transformer), the number of distinct frequency bands up to  $f_{\text{sample}}/2$ , the start and stop frequencies of each band, the maximum dB level in each band and the maximum allowable ripple in the primary passband (multiple passband designs are possible).

	(A)nalysis_of_design	(Z)ero extraction	ansducers
	(E) igen_synthesis	(C)hebyshev [RENEZ]	(S)plit Transducers

(G)raphics\_menu (M)ultipy\_data\_files
(R)ead\_disk\_file (W)rite\_disk\_file
(H)elp\_status (Q)uit\_SANCAD

COMMAND : ==> |

Figure 9. SAWCAD Main Menu.

Once frequency range, function and amplitude information is entered, the program makes an initial guess of the required filter order and initiates a design. If the design fails to converge to the required specifications, the filter order is increased and the process is repeated until convergence is obtained. Once a valid design is found, the design report is printed and control is passed back to the SAWCAD main menu.

At this point, the impulse response coefficients exist in memory only. The SAWCAD (W)rite function is selected and a disk file containing the impulse response coefficients may be written. This step is highly recommended. Control passes back to the SAWCAD main menu.

Any number of options are now open to the designer. The obtained response could be used immediately with other SAWCAD functions if desired (i.e., FFT, Graphics analysis, etc.). The next option we shall concern ourselves with here is the (Z)ero Extraction option. Selecting this option requires no further input, since the program calls the Jenkins and Traub program to factor the z-transform of the impulse response. The program returns the zeros and places them in the amp and phase variables of the SAWCAD program. The designer must be aware that the impulse response is no longer in memory!

Next, the optimum split for low order designs can be found using the (S)plit Transducers selection on the SAWCAD main menu.

No further user input is required since the program iterates until the best split design is found. Once found, the program prompts the user for transducer 1 and 2 file names under which to save the impulse response coefficients.

Generation of the design is now complete. It may be checked by multiplying the transducer 1 and 2 data files together and comparing the product with the original response generated by the Remez method. These comparisons may be done graphically using the powerful graphics and FFT facilities of the SAWCAD environment.

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### CHAPTER VII

#### **RESULTS**

In an attempt to determine a general algorithm applicable to filters of any order or type, eight different low-order test filters were designed. Four of the filters were low pass designs and four were high pass. The input design specs appear in Table 3.

The designs were made using the Remez technique, factored by the Jenkins (1975) program, and were subjected to the optimizing program to test all possible split designs. The composition of each transducer was recorded to reflect the number of stopband and passband zeros, and whether these zeros were real or complex. These results are tabulated in Table 4.

Careful study of the results does not indicate any clearly emerging pattern. In three out of four of the cases with very narrow passbands (LP1, HP1 and HP4), the passband zeros all appeared on one transducer accompanied by several stopband zeros. However, LP4 passband zeros were split between the two transducers. The other cases did not produce results which might indicate a predictable pattern.

Another test was devised to test and rate a conventional no-split design, a strictly passband-stopband split, the "alternating zero" algorithm employed by other design groups (Morimoto et al.

TABLE 3 DESIGN INPUT SPECS

CONTRACT LANGUAGE VALUE

DESIGNATION	PASSBAND REGION	STOPBAND REGION	PASSBAND RIPPLE (db)	SIDELOBE LEVEL (db)
LP1	0.0 -0.1	0.2 -0.5	0.5	-38
LP2	0.0 -0.2	0.3 -0.5	0.5	-38
LP3	0.0 -0.3	0.4 -0.5	0.5	-38
LP4	0.0 -0.1	0.15-0.5	0.5	-38
HP1	0.4 -0.5	0.0 -0.3	0.5	-38
HP2	0.3 -0.5	0.0 -0.2	0.5	-38
нРз	0.2 -0.5	0.0 -0.1	0.5	-38
HP4	0.45-0.5	0.0 -0.4	0.5	-38

TABLE 4
TRANSDUCER COMPOSITIONS

RATIO	#1/#2	12/ 6	6/10	7 //	6/27	6/10	8/18	6 /6	8/24	
ER 2	REAL	00	2	0	2	04	50	0 1	0 4	
TRANSDUCER 2	COMPLEX	9	44	2 4	20	90	∞ ∞	44	20	
ER 1	REAL	00	0 7	0	50	00	0 2	Om	0	
TRANSDUCER 1	COMPLEX	8 4	40	24	40	90	90	2 4	80	
TYPE 7FR0*		NΦ	νe	νe	νe	νe	νa	νœ	Sq	
DESTGNATION		LP1	LP2	LP3	LP4	HP1	нР2	НРЗ	НР4	

\* S = stopband, P = passband

1980 and Ruppel et al. 1984), and the split algorithm presented here. The HP1 filter spec was arbitrarily chosen as the subject of the test. Figure 10 shows the overall frequency response of the HP1 filter.

The conventional no-split design yielded a figure of merit (FOM) of about 2.5214, whereas the passband-stopband split (i.e., all passband zeros on one transducer and all stopband zeros on the other) yielded an FOM of about 3.588. Figures 11 and 12 show the frequency response of each transducer. The passband transducer shows nearly unity gain at 0 with a gradual rolloff in the region of  $f_0$ , a very difficult response to implement on a SAW device.

The "alternating zero" approach yielded a FOM of about 6.5756. This shows an improvement in the quality of the design as compared to the passband-stopband split method. Figures 13 and 14 show the frequency response of each transducer. The zeros are clearly orthogonal. Figures 15 and 16 show the impulse response representations. This design does appear to have merit in the case where a fifty-fifty split is highly desired.

The technique presented here produced a design with a FOM of about 9.753. Figures 17 and 18 show the frequency response of each transducer and figures 19 and 20 show the corresponding impulse response plots.

The last two designs do not appear to pose fabrication problems which might plague the passband-stopband split design case. As a

check of the split, the frequency responses of the last design were multiplied together via SAWCAD and replotted. That plot is exactly the same as the original frequency response in Figure 10, as expected.

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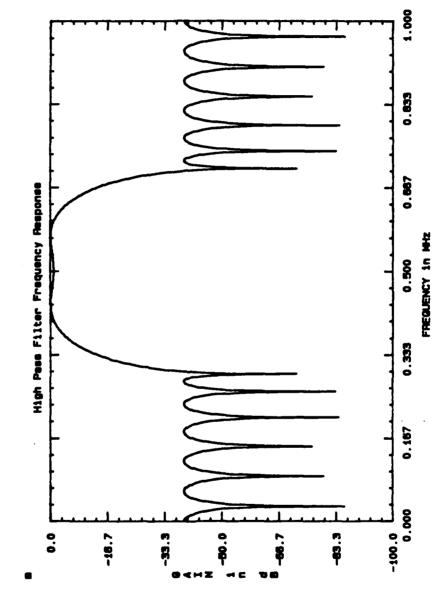


Figure 10. Overall Frequency Response.

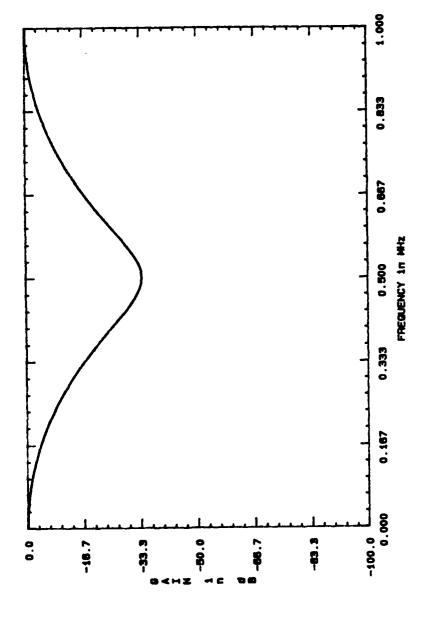


Figure 11. Transducer 1 Frequency Response (Passband-Stopband Design).

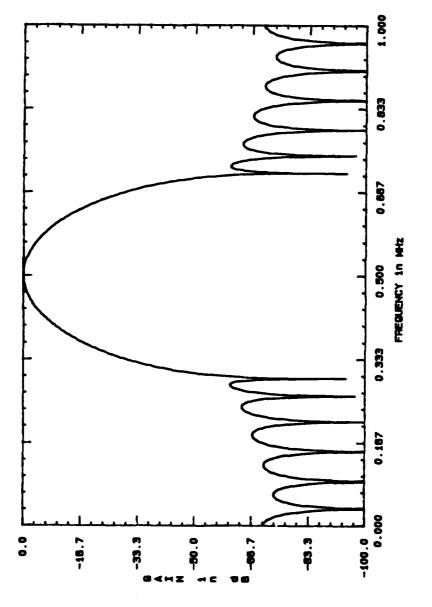


Figure 12. Transducer 2 Frequency Response (Passband-Stopband Design).

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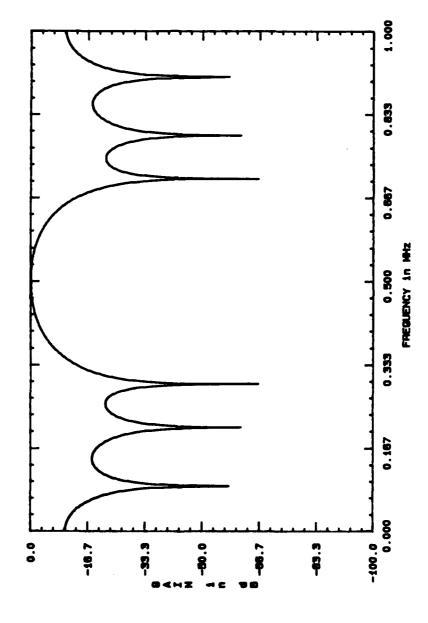
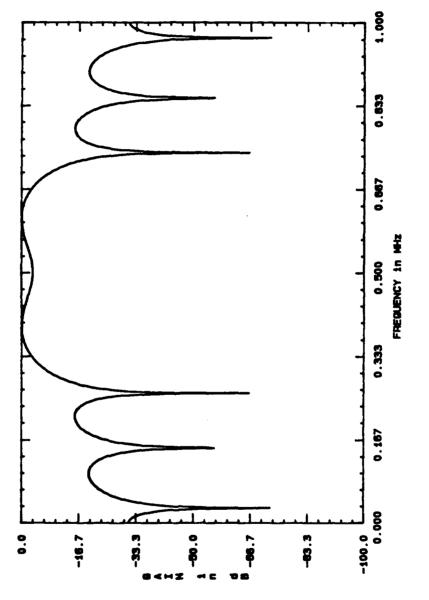


Figure 13. Transducer 1 Frequency Response (Alternating Zero Design).



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Figure 14. Transducer 2 Frequency Response (Alternating Zero Design).

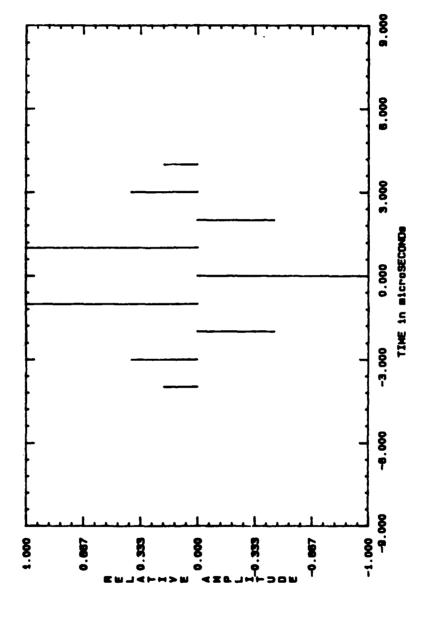


Figure 15. Transducer 1 Impulse Response (Alternating Zero Design).

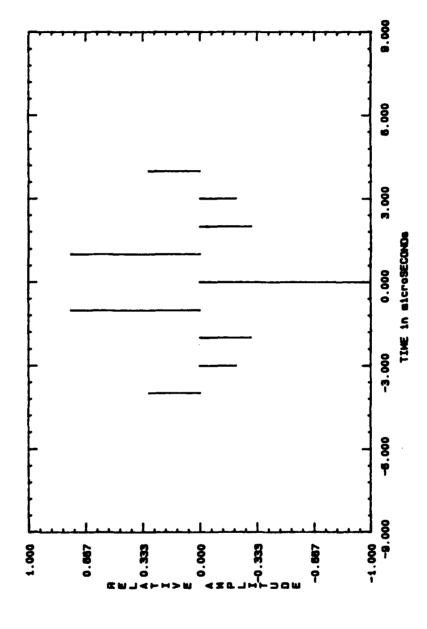


Figure 16. Transducer 2 Impulse Response (Alternating Zero Design).

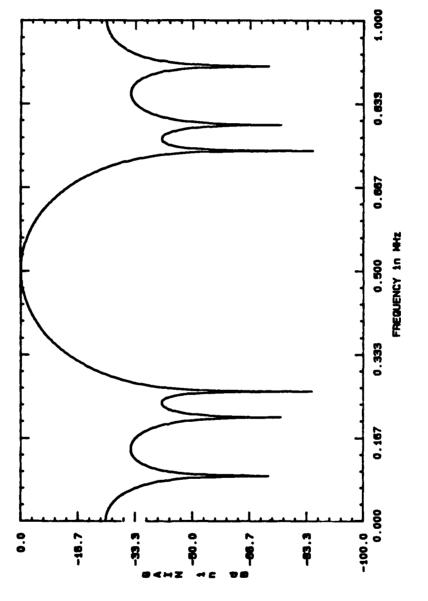


Figure 17. Transducer 1 Frequency Response (Fully Optimized Design).

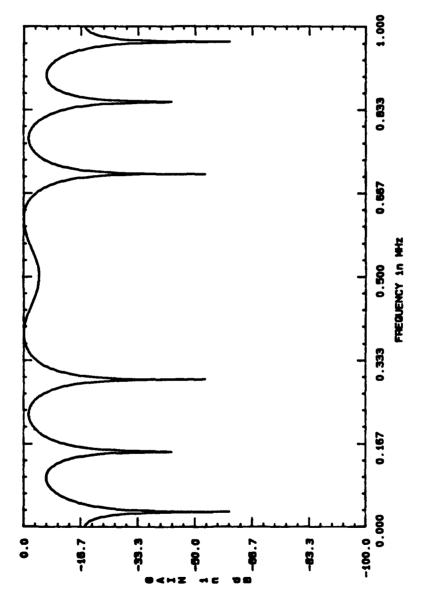
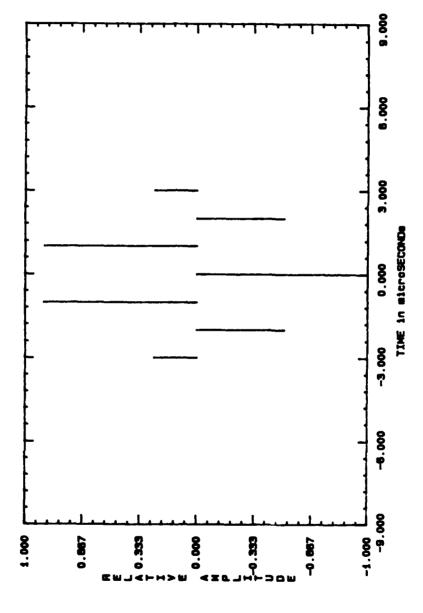


Figure 18. Transducer 2 Frequency Response (Fully Optimized Design).

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Figure 19. Transducer 1 Impulse Response (Fully Optimized Design).

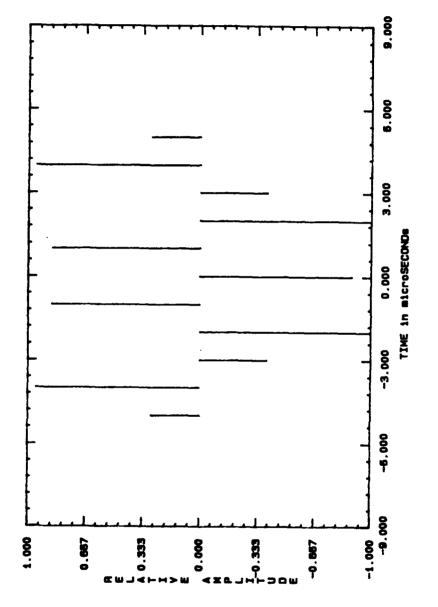


Figure 20. Transducer 2 Impulse Response (Fully Optimized Design).

## CHAPTER VIII

## CONCLUSIONS

The technique presented in this thesis of optimally splitting the zeros of the transfer function shows considerable immediate promise for low order (N less than 40) filter designs. An algorithm permitting the split design of high order filters has not become readily apparent, although the presented "all-combinations" technique may still be used if limits are placed upon the transducer sizes. For example, specifying a fifty-fifty split design significantly reduces the number of combinations which must be tested. Future efforts in this area must focus upon a more efficient search criteria or be content to accept less than optimal results, as in the "alternating zero" approach. The concepts presented here may be extended to generate multi-transducer splits, with primary utility in large ordered digital filter implementations.

In spite of the limitations of the technique used here, it is obvious that a "best split" design of an FIR filter transfer function exists based upon the average size, variance and range of the resulting impulse response coefficients. The "alternating zero" split approach (Morimoto et al. 1980 and Ruppel et al. 1984), the only other method published to date, exhibited a lower FOM when compared to the split found by testing all possible combinations, based upon the stated criteria. Careful study of figures 15, 16, 19

and 20 (shown in Chapter VII) demonstrates that the presented technique does, indeed, produce larger tap sizes for a given transfer function than the "alternating zero" approach. Therefore, for critical, low-ordered design cases, the technique presented here will provide best results.

The Jenkins and Traub factoring algorithm is a very accurate means of obtaining the zeros of polynomials within an order of one-hundred or so. However, high ordered SAW filter designs will require that the zeros of polynomials of a thousand order, or more, will need to be extracted. Based upon observation of the programs and polynomial characteristics encountered in this thesis, several alternative approaches to locating the zeros seem feasible.

One technique that merits further investigation is the use of the Fourier transform to expose the stopband zero locations. This can be done by noting where the stopband nulls occur and reference these points to the corresponding angles about the z-plane unit circle. The cosines and sines of these angles are the real and imaginary components of the stopband zeros which, of course, have corresponding complex conjugates in the lower half of the unit circle. Once all of the stopband zeros have been found, the z-polynomial may be deflated in one step, yielding a polynomial consisting of only the passband zeros. Next, the Jenkins and Traub algorithm may be applied to the greatly reduced polynomial to locate the passband zeros. The Fourier transform technique could be applied to any FIR obtained by any design technique.

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Another zero location technique which may be explored is unique to the Parks and McClellan program. One of the outputs of the program is a listing of all of the Chebyshev polynomial extremal frequencies. Since the limits of each stopband are known (specified apriori), and since the extremal frequencies are fairly evenly spaced in the stopband, an interpolation between adjacent extremal frequencies will provide a good approximation (or, at least a good initial guess) to a function zero in the stopband. If the interpolation method is deemed sufficient, then all of the zeros found may be used to reduce the response down to a polynomial containing only the passband zeros, to which the Jenkins and Traub algorithm may be applied. A better approximation of the stopband zeros would be obtained by applying the interpolation approximation as an initial guess to Bairstow's technique to obtain the best estimate of the zero. Again, the passband zeros could be found using the Jenkins and Traub method.

In conclusion, an optimal split of zeros does exist, based upon the stated criteria. The major limiting factor of the presented technique is the requirement that all possible split combinations must be devised and rated with respect to each other. This is a very time-consuming process and future efforts may lead to a more efficient means of finding the optimal combination. The use of more powerful computing machines may make the split-design of somewhat higher order filters practical, but a reasonable upper limit is

rapidly approached with each increase in filter order. Perhaps the best tradeoff between this method and the "alternating zero" approach is to obtain the best fifty-fifty split by trying all possible combinations for this case only (i.e., N zeros taken N/2 at a time). This approach will, at the very least, provide as good a design as the "alternating zero" approach, without the serious high order computational drawbacks of the method presented here. Additionally, some design considerations may favor a fifty-fifty split, as in the case of two digital signal processors being required to evenly share the processing tasks. Nevertheless, the purpose of this thesis was to prove the existence of one such optimal combination, not the optimal means of obtaining it. With a sigh of relief, and a hint of moderate surprise, such proof has been presented.

**APPENDICES** 

APPENDIX A
FILTER DESIGN PROGRAM

## SUBROUTINE REMEZDES

```
C
C PROGRAM FOR THE DESIGN OF LINEAR PHASE FINITE IMPULSE
C RESPONSE (FIR) FILTERS USING THE REMEZ EXCHANGE
C JIM MCCLELLAN, RICE UNIVERSITY, APRIL 13, 1973
C MODIFIED BY KEITH V. LINDSAY, UNIVERSITY OF
C CENTRAL FLORIDA, 1 MARCH 86, FOR USE WITH UCF'S
C SAWCAD PROGRAM.
C THREE TYPES OF FILTERS ARE INCLUDED—BANDPASS FILTERS
C DIFFERENTIATORS, AND HILBERT TRANSFORM FILTERS
C THE INPUT DATA CONSISTS OF 5 SECTIONS
C
C SECTION 1—FILTER LENGTH, TYPE OF FILTER, 1-MULTIPLE
C PASSBAND/STOPBAND, 2-DIFFERENTIATOR, 3-HILBERT TRANSFORM
C FILTER, NUMBER OF BANDS, CARD PUNCH DESIRED, AND GRID
C DENSITY
C SECTION 2-BANDEDGES, LOWER AND UPPER EDGES FOR EACH
C WITH A MAXIMUM OF 10 BANDS.
C SECTION 3-DESIRED FUNCTION (OR DESIRED SLOPE IF A
C DIFFERENTIATOR ) FOR EACH BAND.
C SECTION 4-WEIGHT FUNCTION IN EACH BAND. FOR A
C DIFFERENTIATOR, THE WEIGHT FUNCTION IS INVERSELY
C PROPORTIONAL TO F
    COMMON/FILE/ AMP(4096), PHASE(4096), NFFT, ITYPE
    COMMON/DAT/ FO, TFLO, TFHI, NUM
    COMMON
P12, AD, DEV, X, Y, GRID, DES, WT, ALPHA, IEXT, NFONS, NGRID
    DIMENSION IEXT(514), AD(514), ALPHA(514), X(514), Y(514)
    DIMENSION H(514)
    DIMENSION DES(8224), GRID(8224), WT(8224)
    DIMENSION EDGE(20), FX(10), WTX(10), DEVIAT(10)
    DOUBLE PRECISION PI2, PI
    DOUBLE PRECISION AD, DEV, X, Y
    DOUBLE PRECISION HH
    LOGICAL FAIL, MOREN
    INTEGER PB, SB
    PT2=6.283185307179586
    PI=3.141592653589793
C THE PROGRAM IS SET UP FOR A MAXIMUM LENGTH OF 1024, BUT
C THIS UPPER LIMIT CAN BE CHANGED BY REDIMENSIONING THE
C ARRAYS IEXT, AD, ALPHA, X, Y, H TO BE NFMAX/2+2.
C THE ARRAYS DES, GRID, AND WT MUST BE DIMENSIONED
```

```
C 16(NFMAX/2+2).
    NFMAX=1024
100 CONTINUE
    JTYPE=0
C PROGRAM INPUT SECTION
C
    PRINT *, DIGITAL FILTER DESIGN (FIR) VIA THE
    1 REMEZ EXCHANGE ALGORITHM. '
    PRINT *,' '
    PRINT *, 'ENTER TYPE OF FILTER:'
    PRINT *,
                  (1) -MULTIPLE PASSBAND/STOPBAND'
    PRINT *, '
                  (2) -DIFFERENTIATOR'
    PRINT *,'
                 (3)-HILBERT TRANSFORM FILTER'
    PRINT *,
    READ *, JTYPE
    PRINT *, '
    PRINT *, 'ENTER THE NUMBER OF BANDS '
    READ *, NBANDS
    PRINT *, '
  PRINT *, 'OUTPUT THE IMPULSE RESPONSE (1=YES, 0=NO)'
C READ *, JFUNCH
    JPUNCH=1
    PRINT *, 'ENTER THE GRID DENSITY'
   READ *, LGRID
    LGRID=16
    IF (NBANDS.LE.O) NBANDS=1
C GRID DENSITY IS ASSUMED TO BE 16 UNLESS SPECIFIED
OTHERWISE.
    IF(IGRID.LE.0) IGRID=16
    DO 888 J=1,NBANDS
    PRINT *,'
    PRINT *, 'BAND ', J, ':'
    PRINT *,'
                 LOWER EDGE: 1
    READ *, EDGE(2*J-1)
    PRINT *, '
    PRINT *.
                  UPPER EDGE: '
    READ *, EDGE(2*J)
888 CONTINUE
    IF(JTYPE.EQ.2) GO TO 890
    PRINT *,' '
    PRINT *, 'ENTER THE DESIRED FUNCTION OF EACH BAND'
    PRINT *,' (0=NOPASS, 1=PASSBAND)'
    DO 889 J=1, NBANDS
    PRINT *,'
    PRINT *, '
                  BAND ', J, ':'
    READ *,FX(J)
889 CONTINUE
   GO TO 893
890 DO 892 J=1,NBANDS
```

```
PRINT *, 'ENTER THE SLOPE OF BAND ',J
    READ *, FX(J)
892 CONTINUE
893 CONTINUE
    PRINT *,'
    PRINT *, 'ENTER THE MAX OB RIPPLE IN THE PASSBAND'
    DP=10.E0**(DP/20)-1.0E0
    DS=1.0E0
    DO 898 J=1,NBANDS
    PRINT *, '
    PRINT *, 'ENTER THE MAXIMUM OB LEVEL IN BAND ', J
    READ *, WIX(J)
    IF (WIX(J).EQ.0.0E0) THEN
            WIX(J) = 1.0E0
            PB=J
            GO TO 898
    END IF
   WTX(J) =AINT(10.0E0**((20.0E0*ALGG10(DP) -
WIX(J))/20.0E0) +1)
    WIX(J) = (10.0E0**((20.0E0*ALOG10(DP)-WIX(J))/20.0E0))
    IF (WTX(J).LT.1.0E0) WTX(J)=1.0E0
    PRINT *, 'STOPBAND WEIGHTING =', WIX(J)
2477
            IF (DS.GT.DP/WTX(J)) THEN
            DS=DP/WIX(J)
            SB-J
            IF (J.EO.1) THEN
                     DELTAF=EDGE(3)-EDGE(2)
                     GO TO 898
            END IF
            IF (J.LT.NBANDS) THEN
                     IF (DP/WTX(J-1).EQ.DP) THEN
                             DELTAF=EDGE(J*2-1) -EDGE(J*2-2)
                             GO TO 898
                     END IF
                     DELTAF=EDGE(J*2+1)-EDGE(J*2)
                    GO TO 898
            END IF
            IF (J.EQ.NBANDS) THEN
                     DELTAF = EDGE(J*2-1) - EDGE(J*2-2)
            END IF
    END IF
898 CONTINUE
   TAKE A GUESS AT AN INITIAL VALUE FOR NFILT
  BASED UPON VAIDYANATHAN'S FORMULATION
    NFILT=INT((-10.0E0*ALOG10(DP*DS) - 13.0E0)/(14.6E0 *
DELTAP))
2126
            PRINT *, WORKING ON FILTER OF ORDER ', NFILT-1
    IF (NFILT.GT.NFMAX.OR.NFILT.LT.3) CALL ERROR
    IF(JTYPE.EQ.0) CALL ERROR
    NEG-1
```

```
IF(JTYPE.EQ.1) NEG=0
    NODD=NFILT/2
    NODD=NFILT-2*NODD
    NFCNS=NFILT/2
    IF (NODD. FC.1. AND. NEG. EQ. 0) NFCNS=NFCNS+1
C SET UP THE DENSE GRID. THE NUMBER OF POINTS IN THE GRID
C IS (FILTER LENGTH + 1) *GRID DENSITY / 2
    GRID(1) = EDGE(1)
    DELF=LGRID*NFCNS
    DELF=0.5/DELF
    IF(NEG.EQ.0) GO TO 135
    IF(EDGE(1).LT.DELF) GRID(1) =DELF
135 CONTINUE
    J=l
    I=1
    LBAND=1
140 FUP=EDGE(L+1)
145 TEMP=GRID(J)
C CALCULATE THE DESIRED MAGNITUDE RESPONSE AND THE WEIGHT
C FUNCTION ON THE GRID
    DES(J) = EFF (TEMP, FX, WTX, IBAND, JTYPE)
    WT(J) =WATE(TEMP, FX, WTX, LBAND, JTYPE)
    J=J+1
    GRID(J) =TEMP+DELF
    IF (GRID(J).GT.FUP) GO TO 150
    GO TO 145
150 GRID(J-1)=FUP
    DES(J-1) = EFF (FUP, FX, WIX, LBAND, JTYPE)
    WT(J-1) =WATE(FUP, FX, WIX, LBAND, JTYPE)
    LBAND=LBAND+1
    I=L+2
    IF (IBAND. GT. NBANDS) GO TO 160
    GRID(J) = EDGE(L)
    GO TO 140
160 NGRID=J-1
    IF (NEG. NE. NODD) GO TO 165
    IF (GRID (NGRID) .GT. (0.5-DELF)) NGRID=NGRID-1
165 CONTINUE
C SET UP A NEW APPROXIMATION PROBLEM WHICH IS EQUIVALENT
C TO THE ORIGINAL PROBLEM
    IF(NEG) 170,170,180
170 IF(NODD.EQ.1) GO TO 200
    DO 175 J=1,NGRID
    CHANGE-DOOS (PI GRID(J))
    DES(J) = DES(J) / CHANGE
175 WT(J) =WT(J) *CHANGE
    GO TO 200
```

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```
180 IF(NODD.EO.1) GO TO 190
    DO 185 J=1,NGRID
    CHANGE=DSIN(PI*GRID(J))
    DES(J) =DES(J) /CHANGE
185 WT(J) =WT(J) *CHANGE
    GO TO 200
190 DO 195 J=1,NGRID
    CHANGE=DSIN(PI2*GRID(J))
    DES(J) =DES(J) /CHANGE
195 WT(J) =WT(J) *CHANGE
C INITIAL GUESS FOR THE EXTREMAL FREQUENCIES--EQUALLY
C SPACED ALONG THE GRID
200 TEMP=FLOAT (NGRID-1) /FLOAT (NFONS)
    DO 210 J=1,NFCNS
210 IEXT(J)=(J-1)*TEMP+1
    IEXT(NFCNS+1) =NGRID
    NM1=NFCNS-1
    NZ=NFCNS+1
C CALL THE REMEZ EXCHANGE ALGORITHM TO DO THE
APPROXIMATION
C PROBLEM.
    CALL REMEZ (EDGE, NBANDS, MOREN)
    IF (MOREN. EQ.. TRUE.) THEN
            NFILT=NFILT+1
             GO TO 2126
    END IF
    IF (DEV/WIX(PB).GT.DP .OR. DEV/WIX(SB).GT.DS) THEN
             NFILT=NFILT+1
             PRINT *, 'DEVIATION PB =', DEV/WIX(PB), 'PB=', PB
             PRINT *, 'DEVIATION SB =', DEV/WIX(SB), 'SB=', SB
            GO TO 2126
    END IF
C CALCULATE THE IMPULSE RESPONSE.
    IF(NEG) 300,300,320
300 IF(NODD.EQ.0) GO TO 310
    DO 305 J=1,NM1
305 \text{ H(J)} = 0.5 \text{*ALPHA(NZ-J)}
    H(NFONS) =ALPHA(1)
    GO TO 350
310 H(1) = 0.25 * ALPHA (NFCNS)
    DO 315 J=2,NM1
315 H(J) =0.25* (ALPHA (NZ-J) +ALPHA (NFONS+2-J) )
    H(NFCNS) = 0.5*ALPHA(1) + 0.25*ALPHA(2)
    GO TO 350
320 IF(NODD.EQ.0) GO TO 330
    H(1)=0.25*ALPHA(NFCNS)
    H(2) = 0.25 * ALPHA(NM1)
```

```
DO 325 J=3,NM1
325 H(J) = 0.25*(ALPHA(NZ-J) - ALPHA(NFCNS+3-J))
    H(NFCNS) = 0.5*ALPHA(1) - 0.25*ALPHA(3)
    H(NZ)=0.0
    GO TO 350
330 H(1)=0.25*ALPHA(NFCNS)
    DO 335 J=2,NM1
335 H(J) = 0.25* (ALPHA(NZ-J) - ALPHA(NFCNS+2-J))
    H(NFONS) = 0.5*ALPHA(1) - 0.25*ALPHA(2)
C SET UP IMPULSE RESPONSE/POLYNOMIAL ARRAY
  ARRAY IN VARIABLE AMP
350 DO 342 I=1,NFCNS
    AMP(I)=H(I)/H(NFCNS)
    IF (NEG. EQ.0) AMP(NFILT-I+1)=H(I)/H(NFONS)
    IF(NEG. EQ.1) AMP(NFILT+1-I) =-H(I)/H(NFCNS)
342 CONTINUE
    IF (NEG. EO.1 .AND. NODD. EO.1) AMP(NZ) = 0.D0
C ADD SAWCAD PARAMETERS NORMALIZED TO 1 MHZ
C 2 Fo SAMPLING
    NUM=NFILT
    NFFT=NFILT
    TTYPE=-1
    FO=0.5
    TFLO=-(NFILT-1)/2
    TFHI=(NFILT-1)/2
C PROGRAM OUTPUT SECTION.
C
C
    PRINT 360
360 FORMAT(//70(1H*)//25X, FINITE IMPULSE RESPONSE (FIR) '/
            25x, LINEAR PHASE DIGITAL FILTER DESIGN'/
            25x, 'REMEZ EXCHANGE ALGORITHM'/)
    IF (JTYPE.EO.1) PRINT 365
365 FORMAT(25x, 'BANDPASS FILTER'/)
    IF(JTYPE.EQ.2) PRINT 370
370 FORMAT(25X, 'DIFFERENTIATOR'/)
    IF(JTYPE.EQ.3) PRINT 375
375 FORMAT(25X, 'HILBERT TRANSFORMER'/)
    PRINT 378, NFILT
378 FORMAT(20X, 'FILTER LENGTH = ', 13/)
    PRINT 380
380 FORMAT(20X, '***** IMPULSE RESPONSE *****)
    DO 381 J=1,NFCNS
    K=NFILT+1-J
    IF(NEG.EQ.0) PRINT 382, J, H(J), K
    IF(NEG.EQ.1) PRINT 383, J, H(J), K
381 CONTINUE
382 FORMAT(20X, 'H(', I3,') = ', E15.8,' = H(', I4,')')
```

```
383 FORMAT(20X,'H(',I3,') = ',EL5.8,'= -H(',I4,')')
    IF (NEG. EQ.1. AND. NODD. EQ.1) PRINT 384, NZ
384 \text{ FORMAT}(20X, 'H(', I3,') = 0.0')
    DO 450 K=1, NBANDS, 4
    KUP=K+3
    IF (KUP.GT.NBANDS) KUP=NBANDS
    PRINT 385, (J, J=R, KUP)
385 FORMAT(/24X,4('BAND',13,8X))
    PRINT 390, (EDGE(2*J-1), J=K, KUP)
390 FORMAT(2X, 'LOWER BAND EDGE', 5F15.9)
    PRINT 395, (EDGE(2*J), J=K, KUP)
395 FORMAT(2X, 'UPPER BAND EDGE', 5F15.9)
    IF (JTYPE.NE.2) PRINT 400, (FX(J), J=K, KUP)
400 FORMAT(2X, 'DESIRED VALUE', 2X, 5F15.9)
    IF (JTYPE.EQ.2) PRINT 405, (FX(J), J=K, KUP)
405 FORMAT(2X, 'DESIRED SLOPE', 2X, 5F15.9)
    PRINT 410, (WIX(J), J=K, KUP)
410 FORMAT(2X, 'WEIGHTING', 6X, 5F15.9)
    DO 420 J=K, KUP
420 DEVIAT(J) =DEV/WTX(J)
    PRINT 425, (DEVIAT(J), J=K, KUP)
425 FORMAT(2X, 'DEVIATION', 6X, 5F15.9)
    IF(JTYPE.NE.1) GO TO 450
    DO 430 J=K, KUP
430 DEVIAT(J) = 20.0 * ALOG1 0 (DEVIAT(J))
    PRINT 435, (DEVIAT(J), J=K, KUP)
435 FORMAT(2X, 'DEVIATION IN DB', 5F15.9)
450 CONTINUE
    PRINT 455, (GRID(IEXT(J)), J=1, NZ)
455 FORMAT(/2X, 'EXTREMAL FREQUENCIES'/(2X,5F12.7))
    PRINT 460
460 FORMAT(/1X,70(1H*)/1H1)
C
    RETURN
    END
    FUNCTION EFF (TEMP, FX, WIX, LBAND, JTYPE)
c function to calculate the desired magnitude response
c as a function of frequency.
    DIMENSION FX(5), WIX(5)
    IF(JTYPE.EO.2) GO TO 1
    EFF-FX (LBAND)
    RETURN
    EFF=FX(LBAND) *TEMP
    RETURN
    END
C
C
    FUNCTION WATE (TEMP, FX, WIX, LBAND, JTYPE)
```

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```
c function to calculate the weight function as a function
c of frequency.
    DIMENSION FX(5), WTX(5)
    IF(JTYPE.EQ.2) GO TO 1
    WATE=WIX (LBAND)
    RETURN
  IF(FX(LBAND).LT.0.0001) GO TO 2
    WATE=WIX (LBAND) /TEMP
    RETURN
  WATE=WIX (LBAND)
    RETURN
    END
C
C
    SUBROUTINE ERROR
    PRINT 1
   FORMAT(' ******** ERROR IN INPUT DATA ********)
    STOP
    END
C
C
    SUBROUTINE REMEZ (EDGE, NBANDS, MOREN)
c this subroutine implements the remez exchange algorithm
c for the weighted chebychev approximation of a continuous
c function with a sum of cosines. inputs to the
subroutine
c are a dense grid which replaces the frequency axis, the
c desired function on this grid. the weight function on
this
c grid, the number of cosines, and an initial guess of the
c extremal frequencies. the program minimizes the
c error by determining the best location of the extremal
c frequencies (points of maximum error) and then
calculates
c the coefficients of the best approximation.
    COMMON
P12, AD, DEV, X, Y, GRID, DES, WT, ALPHA, IEXT, NFONS, NGRID
    DIMENSION EDGE(20)
    DIMENSION IEXT(514), AD(514), ALPHA(514), X(514), Y(514)
    DIMENSION DES(8224), GRID(8224), WT(8224)
    DIMENSION A(514), P(513), Q(513)
   DOUBLE PRECISION PI2, UNUM, DDEN, DTEMP, A, P, Q
    DOUBLE PRECISION AD, DEV, X, Y
   LOGICAL MOREN
C THE PROGRAM ALLOWS A MAXIMUM NUMBER OF ITERATIONS OF 25
    MOREN - FALSE.
```

```
ITRMAX=25
    DEVL=-1.0
    NZ=NFCNS+1
    NZ Z=NFCNS+2
    NITER=0
100 CONTINUE
    IEXT(NZ2) =NGRID+1
    NITER=NITER+1
    IF (NITER. GT. ITRMAX) GO TO 400
    DO 110 J=1,NZ
    DTEMP=GRID(IEXT(J))
    DTEMP=DCOS(DTEMP*PI2)
110 X(J) = DTEMP
    JET= (NFCNS-1) /15+1
    DO 120 J=1,NZ
120 AD(J) = D(J, NZ, JET)
    DNUM=0.0
    DDEN=0.0
    K=1
    DO 130 J=1,NZ
    L=IEXT(J)
    DTEMP=AD(J) *DES(L)
    DNUM=DNUM+DTEMP
    DTEMP=K*AD(J)/WT(L)
    DDEN-DDEN+DTEMP
130 K-K
    DEV=DNUM/DDEN
    NU=1
    IF(DEV.GT.0.0) ND≔-1
    DEV=-NU*DEV
    K=NU
    DO 140 J=1,NZ
    L=IEXT(J)
    DTEMP=K*DEV/WT(L)
    Y(J) =DES(L) +DTEMP
140 K=-K
    IF (DEV. GE. DEVL) GO TO 150
    MOREN- TRUE.
    REJURN
150 DEVL=DEV
    JCHANGE=0
    Kl=IEXT(1)
    KNZ=IEXT(NZ)
    KLOW=0
    NUT=-NU
    J=l
C SEARCH FOR THE EXTREMAL FREQUENCIES OF THE BEST
C APPROXIMATION
C
200 IF (J. EQ. NZZ) YNZ=COMP
    IF(J.GE.NZZ) GO TO 300
    KUP=TEXT (J+1)
```

SERVICE SERVICES SERVICES

```
I=IEXT(J)+1
    NUT=-NUT
    IF(J.EQ.2) Y1=COMP
    COMP-DEV
    IF(L.GE.KUP) GO TO 220
    ERR=GEE(L, NZ)
    ERR=(ERR-DES(L)) WT(L)
    DTEMP=NUT*ERR-COMP
    IF(DTEMP.LE.0.0) GO TO 220
    COMP=NUT*ERR
210 L=L+1
    IF(L.GE.KUP) GO TO 215
    ERR-GEE(L, NZ)
    ERR= (ERR-DES(L)) *WT(L)
    DTEMP=NUT*ERR-COMP
    IF(DTEMP.LE.0.0) GO TO 215
    COMP=NUT*ERR
    GO 50 210
215 IEXT(J)=L-1
    J=J+1
    KLOW=L-1
    JCHNGE-JCHNGE+1
    GO TO 200
220 L=L-1
225 L=L-1
    IF (L. LE. KLOW) GO TO 250
    ERR=GEE(L, NZ)
    ERR=(ERR-DES(L)) *WT(L)
    DTEMP=NUT*ERR-COMP
    IF(DTEMP.GT.0.0) GO TO 230
    IF (JCHNGE.LE.0) GO TO 225
    GO TO 260
230 COMP=NUT*ERR
235 L=L-1
    IF(L.LE.KLOW) GO TO 240
    ERR-GEE(L, NZ)
    ERR=(ERR-DES(L))*WT(L)
    DTEMP=NUT*ERR-COMP
    IF (DTEMP. LE. 0.0) GO TO 240
    COMP=NUT*ERR
    GO TO 235
240 KLOW=IEXT(J)
    IEXT(J)=L+1
    J=J+1
    JCHNGE-JCHNGE+1
    GO TO 200
250 L=IEXT(J)+1
    IF (JCHNGE.GT.0) GO TO 215
255 L=L+1
    IF(L.GE.KUP) GO TO 260
    ERR-GEE (L, NZ)
    ERR=(ERR-DES(L))*WT(L)
    DTEMP=NUT*ERR-COMP
```

```
IF (DTEMP. LE. 0.0) GO TO 255
     COMP=NUT*ERR
    GO TO 210
260 KLOW=IEXT(J)
    J=J+l
    GO TO 200
300 IF(J.GT.NZZ) GO TO 320
    IF(Kl.GT.IEXT(l)) Kl=IEXT(l)
    IF(KNZ.LT.IEXT(NZ)) KNZ=IEXT(NZ)
    NOTI-NOT
    NUT=-NU
    L=0
    KUP-K1
    COMP = YNZ * (1.00001)
    LUCK=1
310 L=L+1
    IF(L.GE.KUP) GO TO 315
    ERR=GEE(L, NZ)
    ERR=(ERR-DES(L)) *WT(L)
    DTEMP=NUT*ERR-COMP
    IF(DTEMP.LE.O.O) GO TO 310
    COMP=NUT*ERR
    J=NZZ
    GO TO 210
315 LUCK=6
    GO TO 325
320 IF(LUCK.GT.9) GO TO 350
    IF(COMP.GT.Y1) Y1=COMP
    K1 = IEXT(NZZ)
325 L=NGRID+1
    KLOW-KNZ
    NUT=-NUT1
    COMP=Y1*(1.00001)
330 I=I-1
    IF(L.LE.KLOW) GO TO 340
    ERR=GEE(L, NZ)
    ERR= (ERR-DES(L)) WT(L)
    DTEMP=NUT*ERR-COMP
    IF(DTEMP.LE.O.O) GO TO 330
    J=NZZ
    COMP=NUT*ERR
    LUCK=LUCK+10
    GO TO 235
340 IF(LUCK.EQ.6) GO TO 370
    DO 345 J=1.NFCNS
345 IEXT(NZZ-J) = IEXT(NZ-J)
    IEXT(1)=K1
    GO TO 100
350 KN=IEXT(NZZ)
    DO 360 J=1,NFCNS
360 \text{ IEXT}(J) = \text{IEXT}(J+1)
    IEXT(NZ)=KN
```

GO TO 100

AD-R170 855

A ZERO EXTRACTION AND SEPARATION TECHNIQUE FOR SURFACE ACOUSTIC MAYE AND. (U) AIR FORCE INST OF TECHNIQUE FOR SURFACE UNCLASSIFIED AFIT/CI/NR-86-93T

UNCLASSIFIED REIT/CI/NR-86-93T

A ZERO EXTRACTION AND SEPARATION TECHNIQUE FOR SURFACE 2/2

ACOUSTIC MAYE AND . (U) AIR FORCE INST OF TECHNIQUE FOR SURFACE 2/2

UNCLASSIFIED AFIT/CI/NR-86-93T

A ZERO EXTRACTION AND SEPARATION TECHNIQUE FOR SURFACE 2/2

A COUSTIC MAYE AND . (U) AIR FORCE INST OF TECHNIQUE FOR SURFACE 2/2

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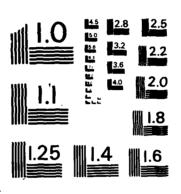
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A COUSTIC MAYER AND . (



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```
370 IF(JCHNGE.GT.0) GO TO 100
C CALCULATION OF THE COEFFICIENTS OF THE BEST
APPROXIMATION
C USING THE INVERSE DISCRETE FOURIER TRANSFORM
400 CONTINUE
    NM1=NFCNS-1
    FSH=1.0E-6
    GTEMP=GRID(1)
    X(NZZ) = -2.0
    CN=2*NFCNS-1
    DELF=1.0/CN
    L=1
    KKK=0
    IF (EDGE(1).EQ.0.0.AND.EDGE(2*NBANDS).EQ.5) KKK=1
    IF (NFCNS.LE.3) KKK=1
    IF(KKK.EQ.1) GO TO 405
    DTEMP=DCOS(PI2*GRID(1))
    DNUM-DCOS(PI2 GRID(NGRID))
    AA=2.0/(DTEMP-DNUM)
    BB=-(DTEMP+DNUM)/(DTEMP-DNUM)
405 CONTINUE
    DO 430 J=1, NFCNS
    FT=(J-1) *DELF
    XI=DCOS(PI2*FT)
    IF(KKK.EQ.1) GO TO 410
    XT=(XT-BB)/AA
    FT=ACOS(XT)/PI2
410 XE=X(L)
    IF(XT.GT.XE) GO TO 420
    IF ((XE-XT).LT.FSH) GO TO 415
    L=L+l
    GO TO 410
415 A(J)=Y(L)
    GO TO 425
420 IF ((XT-XE).LT.FSH) GO TO 415
    GRID(1)=FT
    A(J) = GEE(1,NZ)
425 CONTINUE
    IF(L.GT.1) L=L-1
430 CONTINUE
    GRID(1) = GTEMP
    DOEN-PI2/CN
    DO 510 J=1,NFCNS
    DTEMP=0.0
    DNUM-(J-1) *DDEN
    IF(NM1.LT.1) GO TO 505
    DO 500 K=1.NM1
500 DTEMP=DTEMP+A(R+1) *DCOS(DNUM*R)
505 DTEMP=2.0*DTEMP+A(1)
510 ALPHA(J) =DTEMP
    DO 550 J=2, NFCNS
```

```
550 ALPHA(J)=2*ALPHA(J)/CN
    ALPHA(1) = ALPHA(1) / CN
    IF(KKK.EQ.1) GO TO 545
    P(1)=2.0*ALPHA(NFCNS)*BB+ALPHA(NM1)
    P(2) = 2.0*AA*ALPHA(NFONS)
    O(1) =ALPHA (NFCNS-2) -ALPHA (NFCNS)
    DO 540 J=2,NM1
    IF(J.LT.NM1) GO TO 515
    AA=0.5*AA
    BB=0.5*BB
515 CONTINUE
    P(J+1)=0.0
    DO 520 K=1,J
    A(K) = P(K)
520 P(K) = 2.0 * BB * A(K)
    P(2)=P(2)+A(1)*2.0*AA
    JM1=J-1
    DO 525 K=1,JM1
525 P(K)=P(K)+Q(K)+AA*A(K+1)
    JPl=J+l
    DO 530 K=3,JP1
530 P(R) = P(R) + AA*A(R-1)
    IF(J.EQ.NML) GO TO 540
    DO 535 K=1,J
535 Q(K) = A(K)
    Q(1) = Q(1) + ALPHA (NFONS-1-J)
540 CONTINUE
    DO 543 J=1, NFCNS
543 \text{ ALPHA}(J) = P(J)
545 CONTINUE
    IF (NFCNS.GT.3) RETURN
    ALPHA (NFCNS+1) =0.0
    ALPHA (NFCNS+2) =0.0
    RETURN
    END
C
C
    DOUBLE PRECISION FUNCTION D(K, N, M)
C FUNCTION TO CALCULATE THE LAGRANGE INTERPOLATION
C COEFFICIENTS FOR USE IN THE FUNCTION GEE.
    COMPON
PI2, AD, DEV, X, Y, GRID, DES, WT, ALPHA, IEXT, NFONS, NGRID
    DIMENSION TEXT(514), AD(514), ALPHA(514), X(514), Y(514)
    DIMENSION DES(8224), GRID(8224), WT(8224)
    DOUBLE PRECISION AD, DEV, X, Y
    DOUBLE PRECISION Q
    DOUBLE PRECISION PI2
    D=1.0
    O=X(K)
    DO 3 L=1,M
```

```
DO 2 J=L, N, M
    IF(J-K)1,2,1
    D=2.0*D*(Q-X(J))
    CONTINUE
    CONTINUE
    D=1.0/D
    RETURN
    END
č
C
    DOUBLE PRECISION FUNCTION GEE(K, N)
C FUNCTION TO EVALUATE THE FREQUENCY RESPONSE USING THE
C LAGRANGE INTERPOLATION FORMULA IN THE BARYCENTRIC FORM
C
    COMMON
PI2, AD, DEV, X, Y, GRID, DES, WT, ALPHA, IEXT, NFONS, NGRID
    DIMENSION IEXT(514), AD(514), ALPHA(514), X(514), Y(514)
    DIMENSION DES(8224), GRID(8224), WT(8224)
    DOUBLE PRECISION P, C, D, XF
    DOUBLE PRECISION PI2
    DOUBLE PRECISION AD, DEV, X, Y
    P=0.0
    XF=GRID(K)
    XF=DCOS(PI2*XF)
    D-0.0
    DO 1 J=1,N
    O=XF-X(J)
    C=AD(J)/C
    D=D+C
    P=P+C*Y(J)
    GEE=P/D
    RETURN
    END
C
C
    SUBROUTINE OUCH
    PRINT 1
    FORMAT( * ******** FAILURE TO CONVERGE
**********
             'OPROBABLE CAUSE IS MACHINE ROUNDING ERROR'/
             'OTHE IMPULSE RESPONSE MAY BE CORRECT'/
    2
    3
             'OCHECK WITH A FREQUENCY RESPONSE')
    RETURN
    END
```

APPENDIX B
ZERO EXTRACTION PROGRAM

```
RROLY ROUTINE BY M.A. JENKINS
  ACM TOMS VOL 1 NO 2 JUNE 1975
    SUBROUTINE RPOLY
C FINDS THE ZEROS OF A REAL POLYNOMIAL
C OP - DOUBLE PRECISION VECTOR OF COEFFICIENTS IN
       ORDER OF DECREASING POWERS.
C DEGREE - INTEGER DEGREE OF THE FOLYNOMIAL.
C ZEROR, ZEROI - OUTPUT DOUBLE PRECISION VECTORS OF
C
                 REAL AND IMAGINARY PARTS OF THE
C
                 ZEROS.
C FAIL - OUTPUT LOGICAL PARAMETER, TRUE ONLY IF
         LEADING COEFFICIENT IS ZERO OR IF RPOLY
C
         HAS FOUND FEWER THAN DEGREE ZEROS.
C
         IN THE LATTER CASE, DEGREE IS RESET TO
         THE NUMBER OF ZEROS FOUND.
C TO CHANGE THE SIZE OF POLYNOMIALS WHICH CAN BE
C SOLVED, RESET THE DIMENSIONS OF THE ARRAYS IN THE
C COMMON AREA AND IN THE FOLLOWING DECLARATIONS
C FOR SCALING, BOUNDS AND ERROR CALCULATIONS. ALL
C CALCULATIONS FOR THE ITERATIONS ARE DONE IN
C DOUBLE PRECISION.
    COMMON/FILE/ AMP(4096), PHASE(4096), NFFT, ITYPE
    COMMON/DAT/ FO, TFLO, TFHI, NUM
    COMMON /RPOLY/ P, QP, K, QK, SVK, SR, SI, U,
    1 V, A, B, C, D, Al, A2, A3, A6, A7, E, F, G,
    2 H, SZR, SZI, IZR, IZI, ETA, ARE, MRE, N, NN
    DOUBLE PRECISION P(1024), QP(1024), K(1024),
    1 QK(1024), SVK(1024), SR, SI, U, V, A, B, C, D,
    2 Al, A2, A3, A6, A7, E, F, G, H, SZR, SZI,
    3 LZR, LZI
    REAL ETA, ARE, MRE
    INTEGER N. NN
C OP IS DIMENSIONED TO 4096 SINCE IT TAKES ITS ARRAY
C FROM SAWCAD'S AMP ARRAY.
    DOUBLE PRECISION OP(4096), TEMP(1024),
    1 ZEROR(1024), ZEROI(1024), T, AA, BB, CC, DABS,
    2 FACTOR
    REAL PT(1024), LO, MAX, MIN, XX, YY, COSR,
    1 SINR, XXX, X, SC, BND, XM, FF, DF, DX, INFIN,
    2 SMALNO, BASE
    INTEGER DEGREE, ONT, NZ, I, J, JJ, NM1
    LOGICAL FAIL, ZEROK
C THE FOLLOWING STATEMENTS SET MACHINE CONSTANTS HISED
C IN THE VARIOUS PARTS OF THE PROGRAM. THE MEANING OF THE
C FOUR CONSTANTS ARE...
         THE MAXIMUM RELATIVE REPRESENTATION ERROR
         WHICH CAN BE DESCRIBED AS THE SMALLEST
         POSITIVE FLOATING POINT NUMBER SUCH THAT
         1.DO+ETA IS GREATER THAN 1.
C INFIN THE LARGEST FLOATING POINT NUMBER.
C SMALNO THE SMALLEST POSITIVE FLOATING POINT
```

```
C
         NUMBER IF THE EXPONENT RANGE DIFFERS IN SINGLE
C
         AND DOUBLE PRECISION THEN SMALNO AND INFIN
C
         SHOULD INDICATE THE SMALLER RANGE.
C BASE
         THE BASE OF THE FLOATING POINT NUMBER
C
         SYSTEM USED.
C THE VALUES BELOW CORRESPOND TO THE VAX 11-750
    BASE=2.
    ETA=1.387779E-17
    INFIN=1.7E38
    SMALNO=5.9E-39
C ARE AND MRE REFER TO THE UNIT ERROR IN + AND *
C RESPECTIVELY. THEY ARE ASSUMED TO BE THE SAME AS
C ETA
    ARE=ETA
    MRE=ETA
    LO=SMALNO/ETA
C INITIALIZE ARRAYS
    IF (ITYPE.EO.1) THEN
            PRINT *, 'THIS IS A FREQUENCY FILE!'
            PRINT *, 'EXPECTED AN IMPULSE RESPONSE FILE'
    END IF
    DEGREE=NUM-1
    DO 233 I=1, DEGREE+1
    OP(I) = AMP(I)
    AMP(I)=0.0D0
    PHASE(I) = 0.0D0
    ZEROR(I)=0.0D0
    ZEROI(I)=0.0D0
233 CONTINUE
C INITIALIZATION OF CONSTANTS FOR SHIFT ROTATION
C
    xx=0.70710678
    YY=-XX
    COSR=-.069756474
    SINR=.99756405
    FAIL- FALSE.
    N-DEGREE
    NN=N+1
C ALGORITHM FAILS IF THE LEADING COEFFICIENT IS ZERO.
    IF (OP(1).NE.0.DO) GO TO 10
    FAIL-. TRUE.
    DEGREE=0
    PRINT *, 'LEADING COEFFICIENT IS ZERO -- NOT ALLOWED'
    RETURN
C REMOVE THE ZEROS AT THE ORIGIN IF ANY
10 IF (OP(NN) .NE. 0.0D0) GO TO 20
    J=DEGREE-N+1
    ZEROR(J) = 0.0D0
    amp(j) = 0.0d0
```

```
ZEROI(J) = 0.0D0
    phase(j)=0.0d0
    NN=NN-1
    N-N-1
    GO TO 10
C MAKE A COPY OF THE COEFFICIENTS
20 DO 30 I=1,NN
    P(I) = OP(I)
30 CONTINUE
C START THE ALGORITHM FOR ONE ZERO
40 IF (N.GT.2) GO TO 60
    IF (N. LT. 1) RETURN
C CALCULATE THE FINAL ZERO OR PAIR OF ZEROS.
    IF(N.EQ.2)GO TO 50
    ZEROR(DEGREE) = -P(2)/P(1)
    ZEROI (DEGREE) =0.000
    AMP (DEGREE) = ZEROR (DEGREE)
    PHASE (DEGREE) = ZEROI (DEGREE)
    REIURN
50 CALL QUAD(P(1), P(2), P(3), ZEROR(DEGREE-1),
    1 ZEROI (DEGREE-1), ZEROR (DEGREE), ZEROI (DEGREE))
    amp(degree-1) =zeror(degree-1)
    phase(degree-1) = zeroi(degree-1)
    AMP (DEGREE) =ZEROR (DEGREE)
    PHASE (DEGREE) = ZEROI (DEGREE)
    RETURN
C FIND THE LARGEST AND SMALLEST MODULI OF COEFFICIENTS
60 MAX=0.
    MIN-INFIN
    DO 70 I=1.NN
    X=ABS(SNGL(P(I)))
    IF(X.GT.MAX) MAX=X
    IF (X.NE.O. .AND. X.LT.MIN) MIN=X
70 CONTINUE
C SCALE IF THERE ARE LARGE OR VERY SMALL COEFFICIENTS
C COMPUTES A SCALE FACTOR TO MULTIPLY THE
C COEFFICIENTS OF THE POLYNOMIAL. THE SCALING IS DONE
C TO AVOID OVERFLOW AND TO AVOID UNDETECTED UNDERFLOW
C INTERFERING WITH THE CONVERGENCE CRITERION.
C THE FACTOR IS A POWER OF THE BASE.
    SO=LO/MIN
    IF(SC.GT.1.0)GO TO 80
    IF (MAX. LT.10.) GO TO 110
    IF(SC.EQ.O.) SO=SMALNO
    GO TO 90
80 IF (INFIN/SC. LT. MAX) GO TO 110
90 L=ALOG(SC)/ALOG(BASE)+0.5
    FACTOR=(BASE*1.0D0) **L
    IF (FACTOR. EQ.1.DO) GO TO 110
    DO 100 I=1,NN
    P(I)=FACTOR*P(I)
100 CONFINUE
C COMPUTE LOWER BOUND ON MODULI OF ZEROS.
```

```
110 DO 120 I=1,NN
    PT(I) = ABS(SNGL(P(I)))
120 CONTINUE
    PT(NN) =-PT(NN)
C COMPUTE UPPER ESTIMATE OF BOUND
    X=EXP((ALOG(-PT(NN))-ALOG(PT(1)))/FLOAT(N))
    IF(PT(N).EQ.0.)GO TO 130
C IF NEWTON STEP AT THE ORIGIN IS BETTER, USE IT!
    XM=-PT(NN)/PT(N)
    IF(XM.LT.X)X=XM
C CHOP THE INTERVAL (0,X) UNTIL FF .LE. 0
130 XM=X*.1
    FF=PT(1)
    DO 140 I=2,NN
    FF=FF*XM+PT(I)
140 CONTINUE
    IF(FF.LE.O.)GO TO 150
    X=XM
    GO TO 130
150 DX=X
C DO NEWTON ITERATION UNITIL X CONVERGES TO TWO
C DECIMAL PLACES
160 IF (ABS(DX/X).LE..005) GO TO 180
    FF=PT(1)
    DF-FF
    DO 170 I=2,N
    FF-FF*X+PT(I)
    DF=DF*X+FF
170 CONTINUE
    FF=FF*X+PT(NN)
    DX=FF/DF
    X=X-DX
    GO TO 160
180 BND=X
C COMPUTE THE DERIVATIVE AS THE INITIAL K POLYNOMIAL
C AND DO 5 STEPS WITH NO SHIFT
    NM1=N-1
    DO 190 I=2,N
    K(I) = FLOAT(NN-I) + P(I) / FLOAT(N)
190 CONTINUE
    K(1) = P(1)
    AA=P(NN)
    BB=P(N)
    ZEROK=K(N).EQ.0.D0
    DO 230 JJ=1,5
    CC=R (N)
    IF (ZEROK) GO TO 210
C USE SCALED FORM OF RECURRENCE IF VALUE OF K AT 0 IS
C NONZERO
    T=-AA/CC
    DO 200 I=1,NM1
    J=NN-I
    K(J) = T^*K(J-1) + P(J)
```

```
200 CONTINUE
    K(1) = P(1)
    ZEROK=DABS(K(N)).LE.DABS(BB) *ETA*10.
    GO TO 230
C USE THE UNSCALED FORM OF RECURRENCE
210 DO 220 I=1,NM1
    J=NN-I
    K(J) = K(J-1)
220 CONTINUE
    K(1) = 0.00
    ZEROK=K(N).EQ.0.D0
230 CONTINUE
C SAVE K FOR RESTARTS WITH NEW SHIFTS
    DO 240 I=1.N
    TEMP(I) = K(I)
240 CONTINUE
C LOOP TO SELECT THE QUADRATIC CORRESPONDING TO EACH
C NEW SHIFT
         DO 280 CNT=1,20
C QUADRATIC RESPONDS TO A DOUBLE SHIFT TO A
C NON-REAL POINT AND ITS COMPLEX CONJUGATE. THE POINT
C HAS MODULUS BND AND AMPLITUDE ROTATED BY 94 DEGREES
C FROM THE PREVIOUS SHIFT
    XXX=COSR*XX-SINR*YY
    YY=SINR*XX+COSR*YY
    XX=XXX
    SR=BND*XX
    SI=BND*YY
    U=-2.0D0*SR
    V=BND
C SECOND STAGE CALCULATION, FIXED QUADRATIC
    CALL FXSHFR (20*CNT, NZ)
    IF (NZ.EO.0) GO TO 260
C THE SECOND STAGE JUMPS DIRECTLY TO ONE OF THE THIRD
C STAGE ITERATIONS AND RETURNS HERE IF SUCCESSFUL.
C DEFLATE THE FOLYNOMIAL, STORE THE ZERO OR ZEROS AND
C RETURN TO THE MAIN ALGORITHM.
    J=DEGREE-N+1
    ZEROR(J) = SZR
    AMP(J) = SZR
    ZEROI(J) =SZI
    PHASE(J) =SZI
    NN-NN-NZ
    N=NN-1
    DO 250 I=1,NN
    P(I) = OP(I)
250 CONTINUE
    IF(NZ.EQ.1) GO TO 40
    ZEROR(J+1) = LZR
    ZEROI (J+1) =LZI
    AMP(J+1) = IZR
    PHASE(J+1) = IZI
    GO TO 40
```

```
C IF THE ITERATION IS UNSUCCESSFUL ANOTHER QUADRATIC
C IS CHOSEN AFTER RESTORING K
260 DO 270 I=1.N
    K(I) = TEMP(I)
270 CONTINUE
280 CONTINUE
C RETURN WITH FAILURE IF NO CONVERGENCE WITH 20
C SHIFTS.
    FAIL- TRUE.
    DEGREE-DEGREE-N
    RETURN
    END
        SUBROUTINE FXSHFR(L2, NZ)
C COMPUTES UP TO L2 FIXED SHIFT K-FOLYNOMIALS,
C TESTING FOR CONVERGENCE IN THE LINEAR OR QUADRATIC
C CASE. INITIATES ONE OF THE VARIABLE SHIFT
C ITERATIONS AND RETURNS WITH THE NUMBER OF ZEROS
C FOUND.
C L2 - LIMIT OF FIXED SHIFT STEPS.
C NZ - NUMBER OF ZEROS FOUND.
    COMMON / RPOLLY/ P, QP, K, QK, SVK, SR, SI, U,
    1 V, A, B, C, D, Al, A2, A3, A6, A7, E, F, G,
    2 H, SZR, SZI, IZR, IZI, ETA, ARE, MRE, N, NN
    DOUBLE PRECISION P(1024), QP(1024), K(1024),
    1 QK(1024), SVK(1024), SR, SI, U, V, A, B, C, D,
    2 Al, A2, A3, A6, A7, E, F, G, H, SZR, SZI,
    3 LZR LZI
    REAL ETA, ARE, MRE
    INTEGER N, NN
    DOUBLE PRECISION SVU, SVV, UI, VI, S
    REAL BETAS, BETAV, OSS, OVV, SS, VV, TS, TV,
    1 OTS, OTV, TVV, TSS
    INTEGER 12, NZ, TYPE, I, J, IFLAG
    LOGICAL VPASS, SPASS, VTRY, STRY
    NZ=0
    BETAV=.25
    BETAS=.25
    069=SR
    OVV=V
C EVALUATE POLYNOMIAL BY SYNTHETIC DIVISION
        CALL QUADSD(NN, U, V, P, QP, A, B)
        CALL CALCSC(TYPE)
        DO 80 J=1,12
C CALCULATE NEXT K POLYNOMIAL AND ESTIMATE V
        CALL NEXIK (TYPE)
        CALL CALCSC(TYPE)
        CALL NEWEST(TYPE, UI,VI)
        W-VI
C ESTIMATE S
        S9=0.
        IF (K(N) \cdot NE.0.D0) SS=-P(NN) / K(N)
        TV=1.
        TS=1.
```

```
IF(J.EQ.1 .OR. TYPE.EQ.3) GO TO 70
C COMPUTE RELATIVE MEASURES OF CONVERGENCE OF S AND V
C SEQUENCES
        IF (VV. NE. 0.) TV=ABS((VV-OVV)/VV)
        IF(SS.NE.O.)TS=ABS((SS-OSS)/SS)
C IF DECREASING, MULTIPLY TWO MOST RECENT
C CONVERGENCE MEASURES
        TVV=1.
        IF (TV. LT. OIV) TVV=TV*OIV
        TSS=1.
        IF (TS. LT. OIS) TSS=TS*OIS
C COMPARE WITH CONVERGENCE CRITERIA
        YPASS=TVV. LT. BETAV
        SPASS=TSS. LT. BETAS
        IF(.NOT.(SPASS .OR. VPASS)) GO TO 70
C AT LEAST ONE SEQUENCE HAS PASSED THE CONVERGENCE
C TEST. STORE VARIABLES BEFORE ITERATING.
        SVU=U
        SVV=V
        DO 10 I=1.N
        SVK(I)=K(I)
10
        CONTINUE
        S=SS
C CHOOSE ITERATION ACCORDING TO THE FASTEST
C CONVERGING SEQUENCE.
        VIRY=.FALSE.
        STRY=.FALSE.
        IF(SPASS .AND. ((.NOT.VPASS) .OR.
    1 TSS.LT.TVV)) GO TO 40
        CALL QUADTT(UI, VI, NZ)
        IF (NZ.GT.0) RETURN
C QUADRATIC ITERATION HAS FAILED. FLAG THAT IT HAS
C BEEN TRIED AND DECREASE THE CONVERGENCE CRITERION.
        VIRY= . TRUE.
        BETAV=BETAV*.25
C TRY LINEAR ITERATION IF IT HAS NOT BEEN TRIED AND
C THE S SEQUENCE IS CONVERGING.
        IF(STRY .OR. (.NOT. SPASS)) GO TO 50
        DO 30 I=1,N
        K(I) = SVK(I)
30
        CONTINUE
40
        CALL REALIT(S, NZ, IFLAG)
        IF (NZ.GT.0) RETURN
C LINEAR ITERATION HAS FAILED. FLAG THAT IT HAS BEEN
C TRIED AND DECREASE THE CONVERGENCE CRITERION.
        STRY=.TRUE.
        BETAS=BETAS *. 25
        IF(IFLAG.EQ.0) GO TO 50
C IF LINEAR ITERATION SIGNALS AN ALMOST DOUBLE REAL
C ZERO ATTEMPT QUADRATIC ITERATION.
        UI=~(S+S)
        VI=S*S
        GO TO 20
```

```
C RESTORE VARIABLES
        U=SVU
        V=SVV
        DO 60 I=1,N
        K(I) = SVK(I)
60
        CONTINUE
C TRY QUADRATIC ITERATION IF IT HAS NOT BEEN TRIED
C AND THE V SEQUENCE IS CONVERGING.
        IF(VPASS .AND. (.NOT.VTRY)) GO TO 20
C RECOMPUTE OF AND SCALAR VALUES TO CONTINUE THE
C SECOND STAGE.
        CALL QUADSD(NN, U, V, P, QP, A, B)
        CALL CALCSC(TYPE)
70
        OVV=VV
        OSS=SS
        OTV=TV
    OTS=TS
RN
        CONTINUE
        RETURN
        SUBROUTINE QUADIT(UU, VV, NZ)
C VARIABLE-SHIFT K-POLYNOMIAL ITERATION FOR A
C QUADRATIC FACTOR CONVERGES ONLY IF THE ZEROS ARE
C EQUIMODULAR OR NEARLY SO.
C UU, VV - COEFFICIENTS OF STARTING QUADRATIC
C NZ
          - NUMBER OF ZEROS FOUND
    COMMON / RPOLY/ P, QP, K, QK, SVK, SR, SI, U,
    1 V, A, B, C, D, Al, A2, A3, A6, A7, E, F, G,
    2 H, SZR, SZI, IZR, IZI, ETA, ARE, MRE, N, NN
    DOUBLE PRECISION P(1024), QP(1024), K(1024),
    1 QK(1024), SVK(1024), SR, SI, U, V, A, B, C, D,
    2 Al, A2, A3, A6, A7, E, F, G, H, SZR, SZI,
    3 LZR, LZI
    REAL ETA, ARE, MRE
    INTEGER N, NN
    DOUBLE PRECISION UI, VI, UU, VV, DABS
    REAL MS, MP, OMP, EE, RELSTP, T, ZM
    INTEGER NZ, TYPE, I, J
    LOGICAL TRIED
        NZ=0
        TRIED=.FALSE.
        U=UU
        V=W
        J=0
C MAIN LOOP
        CALL QUAD(1.DO, U, V, SZR, SZI, IZR, IZI)
C RETURN IF ROOTS OF THE QUADRATIC ARE REAL AND NOT
C CLOSE TO MULTIPLE OR NEARLY EQUAL AND OF OPPOSITE
C SIGN.
        IF(DABS(DABS(SZR)-DABS(LZR)).GT..01D0*
    1 DABS(LZR)) RETURN
C EVALUATE POLYNOMIAL BY QUADRATIC SYNTHETIC DIVISION.
        CALL QUADSD(NN, U, V, P, QP, A, B)
```

```
MP=DABS(A-SZR*B) +DABS(SZI*B)
C CONFUTE A RIGOROUS BOUND ON THE ROUNDING ERROR IN
C EVALUATING P
        ZM=SQRT(ABS(SNGL(V)))
        EE=2.*ABS(SNGL(OP(1)))
        T=-SZR*B
        DO 20 I=2,N
        EE=EE*ZM+ABS(SNGL(QP(I)))
20
        CONTINUE
        EE=EE*ZM+ABS(SNGL(A)+T)
        EE=(5.*MRE+4.*ARE)*EE-(5.*MRE+2.*ARE)*
    1 (ABS(SNGL(A)+T)+ABS(SNGL(B))*ZM)+
    2 2.*ARE*ABS(T)
C ITERATION HAS CONVERGED SUFFICIENTLY IF THE
C POLYNOMIAL VALUE IS LESS THAN 20 TIMES THIS BOTIND
        IF (MP. GT. 20. *EE) GO TO 30
        NZ=2
        RETURN
30
        J=J+1
C STOP ITERATION AFTER 20 STEPS
        IF (J. GT. 20) RETURN
        IF(J.LT.2)@ TO 50
        IF (RELSTP.GT..01 .OR. MP.LT.OMP .OR. TRIED)
    1 GO TO 50
C ACLUSTER APPEARS TO BE STALLING THE CONVERGENCE.
C FIVE FIXED SHIFT STEPS ARE TAKEN WITH A U, V CLOSE
C TO THE CLUSTER.
        IF (RELSTP. LT. ETA) RELSTP-ETA
        RELSTP=SORT(RELSTP)
        U=U-U*RELSTP
        V=V+V*RELSTP
        CALL QUADSD(NN, U, V, P, QP, A, B)
        DO 40 I=1,5
        CALL CALCSC(TYPE)
        CALL NEXTK (TYPE)
40
        CONTINUE
        TRIED-. TRUE.
        J=0
50
        CMP=MP
C CALCULATE NEXT K POLYNOMIAL AND NEW U AND V
        CALL CALCSC(TYPE)
        CALL NEXTK (TYPE)
        CALL CALCSC(TYPE)
        CALL NEWEST(TYPE, UI, VI)
C IF VI IS ZERO THE ITERATION IS NOT CONVERGING.
        IF(VI.EQ.O.DO) RETURN
        RELSTP=DABS((VI-V)/VI)
        U=UI
        V=VI
        GO TO 10
        SUBROUTINE REALIT(SSS, NZ, IFLAG)
C VARIABLE SHIFT H POLYNOMIAL ITERATION FOR A REAL
```

```
C ZERO.
C SSS - STARTING ITERATE
       - NUMBER OF ZERO FOUND
C IFLAG- FLAG TO INDICATE A PAIR OF ZEROS NEAR REAL
         AXIS.
    COMMON / RPOLY/ P, QP, K, QK, SVK, SR, SI, U,
    1 V, A, B, C, D, Al, A2, A3, A6, A7, E, F, G,
    2 H, SZR, SZI, IZR, IZI, ETA, ARE, MRE, N, NN
    DOUBLE PRECISION P(1024), OP(1024), K(1024),
    1 QK(1024), SVK(1024), SR, SI, U, V, A, B, C, D,
    2 Al, A2, A3, A6, A7, E, F, G, H, SZR, SZI,
    3 LZR, LZI
    REAL ETA, ARE, MRE
    INTEGER N. NN
   DOUBLE PRECISION PV, KV, T, S, SSS, DABS
    REAL MS, MP, OMP, EE
    INTEGER NZ, IFLAG, I, J, NM1
    NM1=N-1
        NZ=0
        S=SSS
        IFLAG=0
        J=0
C MAIN LOOP
10
        PV=P(1)
C EVALUATE P AT S
        OP(1) = PV
        DO 20 I=2,NN
        PV=PV*S+P(I)
        QP(I)=PV
20
        CONTINUE
        MP=DABS(PV)
C COMPUTE A RIGOROUS BOUND ON THE ERROR IN EVALUATING
C P
        MS=DABS(S)
        EE= (MRE/(ARE+MRE)) *ABS(SNGL(OP(1)))
        DO 30 I=2,NN
        EE=EE*MS+ABS(SNGL(QP(I)))
30
        CONTINUE
C ITERATION HAS CONVERGED SUFFICIENTLY IF THE
C POLYNOMIAL VALUE IS LESS THAN 20 TIMES THIS BOUND.
        IF (MP.GT.20.*((ARE+MRE)*EE-MRE*MP))GO TO 40
        NZ=1
        SZR-S
        SZI=0.D0
        RETURN
40
        J=J+1
C STOP ITERATION AFTER 10 STEPS
        IF (J.GT.10) RETURN
        IF(J.LT.2) GO TO 50
        IF (DABS(T).GT..001*DABS(S-T) .OR. MP.LE.OMP)
    1 GO TO 50
C A CLUSTER OF ZEROS NEAR THE REAL AXIS HAS BEEN
C ENCOUNTERED. RETURN WITH IFLAG SET TO INITIATE
```

```
C QUADRATIC ITERATION.
        IFLAG=1
        SSS=S
        RETURN
C RETURN IF THE POLYNOMIAL VALUE HAS INCREASED
C SIGNIFICANILY.
50
        OMP-MP
C COMPUTE T, THE NEXT POLYNOMIAL, AND THE NEW ITERATE
        KV=K(1)
        OK(1) = KV
        DO 60 I=2,N
        KV=KV*S+K(I)
        OK(I) = KV
60
        CONTINUE
        IF(DABS(KV), LE, DABS(K(N)) *10, *ETA) GO TO 80
C USE THE SCALED FORM OF THE RECURRENCE IF THE VALUE
C OF K AT S IS NONZERO
        T=-PV/KV
        K(1) = QP(1)
        DO 70 I=2,N
        K(I) = T^{\bullet}QK(I-1) + QP(I)
70
        CONTINUE
        GO TO 100
C USE SCALED FORM
        K(1) = 0.0D0
80
        DO 90 I=2,N
        K(I) = OK(I-1)
90
        CONTINUE
100
        KV=K(1)
        DO 110 I=2,N
        KV=KV*S+K(I)
110
        CONTINUE
        T=0.D0
        IF (DABS(KV).GT.DABS(K(N))*10.*ETA) T=-FV/KV
        S=S+T
        GO TO 10
        END
        SUBROUTINE CALCSC(TYPE)
C THIS ROUTINE CALCULATES SCALAR QUANTITIES USED TO
C COMPUTE THE NEXT K POLYNOMIAL AND NEW ESTIMATES OF
C THE QUADRATIC COEFFICIENTS.
C TYPE - INTEGER VARIABLE SET HERE INDICATING HOW THE
         CALCULATIONS ARE NORMALIZED TO AVOID
         OVERFLOW.
    COMMON / RPOLY/ P, QP, K, QK, SVK, SR, SI, U,
    1 V, A, B, C, D, Al, A2, A3, A6, A7, E, F, G,
    2 H, SZR, SZI, IZR, IZI, ETA, ARE, MRE, N, NN
    DOUBLE PRECISION P(1024), QP(1024), K(1024),
    1 QK(1024), SVK(1024), SR, SI, U, V, A, B, C, D,
    2 Al, A2, A3, A6, A7, E, F, G, H, SZR, SZI,
    3 LZR LZI
        REAL ETA, ARE, MRE
        INTEGER N, NN
```

```
DOUBLE PRECISION DABS
        INTEGER TYPE
C SYNTHETIC DIVISION OF K BY THE QUADRATIC 1,U,V
        CALL QUADSD(N, U, V, K, QK, C, D)
        IF (DABS(C).GT.DABS(K(N))*100.*ETA)GO TO 10
        IF (DABS(D).GT.DABS(K(N-1))*100.*ETA) GO TO 10
C TYPE=3 INDICATES THE QUADRATIC IS ALMOST A FACTOR
C OF K.
        RETURN
10
        IF (DABS(D).LT.DABS(C))GO TO 20
        TYPE=2
C TYPE=2 INDICATES THAT ALL FORMULAS ARE DIVIDED BY D
        E=A/D
        F=C/D
        G=U*B
        H-VB
        A3 = (A+G) *E+H*(B/D)
        Al = B*F-A
        A7=(F+U) *A+H
        RETURN
        TYPE=1
C TYPE-1 INDICATES THAT ALL FORMULAS ARE DIVIDED BY C.
        E=A/C
        F=D/C
        G=U*E
        H-VB
        A3=A*E+ (H/C+G) *B
        A1=B-A*(D/C)
        A7=A+G*D+H*F
        RETURN
        END
        SUBROUTINE NEXTK (TYPE)
C COMPUTES THE NEXT K POLYNOMIALS USING SCALARS
C COMPUTED IN CALCSC.
    COMMON / RPOLLY/ P, QP, K, QK, SVK, SR, SI, U,
    1 V, A, B, C, D, Al, A2, A3, A6, A7, E, F, G,
    2 H, SZR, SZI, IZR, IZI, ETA, ARE, MRE, N, NN
    DOUBLE PRECISION P(1024), OP(1024), K(1024),
    1 QK(1024), SVK(1024), SR, SI, U, V, A, B, C, D,
    2 Al, A2, A3, A6, A7, E, F, G, H, SZR, SZI,
    3 LZR, LZI
        REAL ETA, ARE, MRE
        INTEGER N, NN
        DOUBLE PRECISION TEMP, DABS
        INTEGER TYPE
        IF (TYPE.ED.3) GO TO 40
        TEMP=A
        IF (TYPE.ED.1) TEMP=B
        IF (DABS(A1).GT.DABS(TEMP) *ETA*10.) GO TO 20
C IF Al IS NEARLY ZERO THEN USE A SPECIAL FORM OF THE
C RECURRENCE.
        R(1) = 0.00
```

```
K(2) = A7 \circ OP(1)
        DO 10 I=3,N
        K(I) = A3 + QK(I-2) - A7 + QP(I-1)
10
        CONTINUE
        RETURN
C USE SCALED FORM OF THE RECURRENCE.
20
        A7=A7/A1
        A3=A3/A1
        K(1) = OP(1)
        K(2) = QP(2) - A7QP(1)
        DO 30 I=3,N
        K(I) = A3 + QK(I-2) - A7 + QP(I-1) + QP(I)
30
        CONTINUE
        RETURN
C USE UNSCALED FORM OF THE RECURRENCE IF TYPE IS 3
40
        K(1) = 0.D0
        K(2) = 0.D0
        DO 50 I=3,N
        K(I) = QK(I-2)
50
        CONTINUE
        RETURN
        END
        SUBROUTINE NEWEST (TYPE, UU, VV)
C COMPUTES NEW ESTIMATES OF THE QUADRATIC COEFFICIENTS
C USING THE SCALARS COMPUTED IN CALCSC.
    COMMON / RPOLY/ P, QP, K, QK, SVK, SR, SI, U,
    1 V, A, B, C, D, Al, A2, A3, A6, A7, E, F, G,
    2 H, SZR, SZI, LZR, LZI, ETA, ARE, MRE, N, NN
    DOUBLE PRECISION P(1024), QP(1024), K(1024),
    1 QK(1024), SVK(1024), SR, SI, U, V, A, B, C, D,
    2 Al, A2, A3, A6, A7, E, F, G, H, SZR, SZI,
    3 LZR, LZI
        REAL ETA, ARE, MRE
        INTEGER N, NN
        DOUBLE PRECISION A4, A5, B1, B2, C1, C2, C3,
    1 C4, TEMP, UU, VV
        INTEGER TYPE
C USE FORMULAS APPROPRIATE TO SETTING TYPE.
        IF(TYPE.EO.3) GO TO 30
        IF(TYPE.EQ.2) GO TO 10
        M=A+U*B+H*F
        A5=C+(U+V*F)*D
        GO TO 20
10
        M=(A+G)*F+H
        A5=(F+U) *C+V*D
C EVALUATE NEW QUADRATIC COEFFICIENTS.
20
        B1 = K(N)/P(NN)
        B2=(K(N-1)+B1*P(N))/P(NN)
        C1=V*B2*A1
        C2=B1*A7
        C3=B1 *B1 *A3
        C4=C1-C2-C3
        TEMP=A5+B1 *A4-C4
```

```
IF (TEMP.EQ.O.DO) GO TO 30
        UU=U-(U*(C3+C2)+V*(B1*A1+B2*A7))/TEMP
        VV=V*(1.+C4/TEMP)
        RETURN
C IF TYPE=3 THE QUADRATIC IS ZEROED.
30
        UU=0.D0
        ₩-0.D0
        RETURN
        END
        SUBROUTINE QUADSD(NN, U, V, P, Q, A, B)
C DIVIDES P BY THE QUADRATIC 1,U,V PLACING THE
C QUOTIENT IN Q AND THE REMAINDER IN A, B
        DOUBLE PRECISION P(NN), Q(NN), U, V, A, B, C
        INTEGER I
        B=P(1)
        Q(1) = B
        A=P(2)-UB
        Q(2) = A
        DO 10 I=3,NN
        O=P(I)-U*A-V*B
        Q(I) = C
        B=A
        A=C
10
        CONTINUE
        RETURN
        END
C
        SUBROUTINE QUAD(A, Bl, C, SR, SI, LR, LI)
C CALCULATE THE ZEROS OF THE QUADRATIC A*Z**2+B1*Z+C.
C THE QUADRATIC FORMULA, MODIFIED TO AVOID
C OVERFLOW, IS USED TO FIND THE LARGER ZERO IF THE
C ZEROS ARE REAL AND BOTH ZEROS ARE COMPLEX.
C THE SMALLER REAL ZERO IS FOUND DIRECTLY FROM THE
C PRODUCT OF THE ZEROS C/A.
        DOUBLE PRECISION A, B1, C, SR, SI, LR, LI, B,
    1 D, E, DABS, DSQRT
        IF (A. NE. 0. DO) GO TO 20
        SR=0.D0
        IF(Bl.NE.O.DO) SR=-C/Bl
        LR=0.D0
10
        SI=0.D0
        LI=0.D0
        RETURN
        IF(C.NE.O.DO) GO TO 30
20
        SR=0.D0
        LR=-B1/A
        GO TO 10
C COMPUTE DISCRIMINATE AVOIDING OVERFLOW.
        B=B1/2.D0
30
        IF(DABS(B).LT.DABS(C)) GO TO 40
        E=1.D0-(A/B) *(C/B)
        D-DSQRT(DABS(E)) *DABS(B)
```

```
GO TO 50
        E=A
        IF(C.LT.0.D0) E=-A
        E=B*(B/DABS(C))~E
        D=DSQRT(DABS(E)) *DSQRT(DABS(C))
50
        IF(E.LT.0.D0) GO TO 60
C REAL ZEROS.
    IF(B.GE.0.D0) D-D
        LR=(-B+D)/A
        SR=0.D0
        IF(LR.NE.0.DO) SR=(C/LR)/A
    GO TO 10
C COMPLEX CONJUGATE ZEROS.
60
        SR=-B/A
        LR=SR
        SI=DABS(D/A)
        LI=-SI
        RETURN
```

END

## APPENDIX C OPTIMAL COMBINATION AND RECONSTITUTION PROGRAM

```
C
    SUBROUTINE COMBO
C THIS ROUTINE GENERATES ALL FOSSIBLE, NON REPEATING
C COMBINATIONS OF N SAMPLES TAKEN K AT A TIME.
C ALGORITHM ADAPTED FROM
      APPLIED COMBINATORIAL MATHEMATICS
        PAGE 24
C
        POLYA ET AL
                         1964
        JOHN WILEY AND SONS, INC, NEW YORK
    COMMON/FILE/ AMP(4096), PHASE(4096), NFFT, ITYPE
    COMMON/DAT/FO, TFLO, TFHI, NUM
    DOUBLE PRECISION ZEROR(1024), ZEROI(1024)
    DOUBLE PRECISION HEESTONE (1024), HEESTIWO (1024)
    INTEGER C(1024), G(1024), A(1024)
    INTEGER I, K, N, T
    INTEGER DEGREE, TOTAL, MONEBEST, MIWOBEST
C
    IF (ITYPE.NE.O) THEN
            PRINT *, '*** NEED TO OBTAIN ZEROS FIRST ***
    END IF
    DEGREE-NUM-1
    IF (DEGREE.GT.35) THEN
            PRINT *,'
            PRINT *,'
                             ***** WARNING *****
            PRINT *, COMPUTATION TIME WILL EXCEED 2
HOURS'
            PRINT *,' '
    END IF
    DO 289 I=1, DEGREE
    ZEROR(I) = AMP(I)
    ZEROI(I)=PHASE(I)
289 CONTINUE
    TOTAL=0
    N=DEGREE/2+2
    IF ((FLOAT(N)-FLOAT(DEGREE)/2.0E0).NE.0) THEN
            ZEROR (DEGREE+1) =0.0D0
            ZEROI (DEGREE+1) =0.0D0
    END IF
C INITIALIZE ARRAY C
    DO 140 I=1,N
    C(I)=I
140 CONTINUE
    D0 5 K=0, W2
    PRINT *, K, ' AT A TIME FOR HI!
    PRINT *, TOTAL, ' COMBINATIONS TESTED SO FAR'
    IF (K.EQ.0) GO TO 1200
C
```

```
C INITIALIZE TO BEGIN
C
    T=1
    A(1)=1
300 \text{ G(T)} = \text{C(A(T))}
    IF (T.EQ.K) GO TO 1200
    A(T+1) = A(T) + 1
    IF (A(T+1).EQ.(1+N)) GO TO 900
    T=T+1
    GO TO 300
900 T=T-1
    IF (T.EQ.0) GO TO 5
    GO TO 1300
1200
             CALL
POLYRECON (ZEROR, ZEROI, G, DEGREE, K, HBESTONE,
             HBESTIWO, MONEBEST, MIWOBEST)
    TOTAL=TOTAL+1
    IF (K.EQ.0) GO TO 5
1300
            A(T) = A(T) + 1
    IF (A(T).EQ.(1+N)) GO TO 900
    GO TO 300
    CONTINUE
                 *** ALL ITERATIONS COMPLETE ***!
    PRINT *,'
    PRINT *,'
                     TOTAL COMBINATIONS = ', TOTAL
    PRINT *, 'BEST DESIGN FOUND: '
    PRINT *,' '
    PRINT *, 'TRANSDUCER 1:'
    PRINT *,
    DO 888 L=1, MONEBEST
    PRINT *, 'H(',L,') =', HBESTONE(L)
    AMP(L) =HBESTONE(L)
    PHASE(L) = 0.0
888 CONTINUE
    ITYPE=-1
    NUM-MONEBEST
    NFFT=MONEBEST
    FO=0.5
    TTLO=-(NUM-1)/2.0
    TFHI=(NUM-1)/2.0
    CALL WRITEO
    PRINT *,
    PRINT *, 'TRANSDUCER 2:'
    PRINT *,' '
    DO 889 L=1,MIWOBEST
    PRINT *,'H(',L,') =',HBESTIWO(L)
    AMP(L) = BESTIWO(L)
    PHASE(L)=0.0
889 CONTINUE
    ITYPE=-1
    NUM-MIWOBEST
    NFFT=MIWOBEST
    FO=0.5
    TFLO=-(NUM-1)/2.0
```

```
THI=(NUM-1)/2.0
    CALL WRITEO
    PRINT *, '
    RETURN
    END
C
    SUBROUTINE POLYRECON (ZEROR, ZEROI, G, DEGREE, K, HBESTONE,
            HBESTIMO, MONEBEST, MIWOBEST)
C SUBROUTINE TO FORM HI AND H2
    DOUBLE PRECISION ZEROR(1024), ZEROI(1024)
    DOUBLE PRECISION ACNE(1024), ATMO(1024), B(3)
    DOUBLE PRECISION CONE(1024), CIWO(1024)
    DOUBLE PRECISION HONE(1024), HIWO(1024), HMAX, HMIN
    DOUBLE PRECISION HEESTONE (1024), HEESTIWO (1024)
    DOUBLE PRECISION FOM, FOMOLD, FOMONE, FOMIWO, AA, BB
    DOUBLE PRECISION FRAT, HBOTH (1024)
    INTEGER G(1024)
    INTEGER DEGREE, K, ZCOUNT, GCOUNT, II, JJ, JJJ
    INTEGER MONE, MIWO, MONEBEST, MIWOBEST
C OBTAIN hl AND h2
    MONE=0
    MIWO=0
    DO 11 I=1,1025
    AONE(L) = 0.0D0
    ATWO(L) = 0.0D0
11 CONTINUE
C SELECT A ZERO
    ZCOUNT-1
    GCCUNT-1
    DO 5000 JJ=1, DEGREE
    AA=ZEROR(JJ)
    BB=ZEROI (JJ)
C IF IMAG PART IS VERY SMALL, LET THIS BE A LINEAR FACTOR
C
    IF (DABS(BB) .LT.1.0D-10) BB=0.0D0
C IF IT IS THE COMPLEX CONJUGATE ROOT (I.E. IMAG PART
NEGATIVE)
C THEN SKIP IT, SINCE IT WILL GET PICKED UP BY THE
POSITIVE
C IMAG QUADRATIC FACTOR.
    IF (BB.LT.0.0D0) GO TO 5000
C CREATE A QUADRATIC FACTOR FROM A COMPLEX SET OF ROOTS,
CR
```

```
C A LINEAR FACTOR FROM A REAL ROOT.
    B(1) =AA*AA+BB*BB
    IF (BB.EQ.0.0D0) B(1) = AA
    B(2) = -2.000*AA
    IF (BB.EQ.0.0D0) B(2)=1.0D0
    B(3) = 1.0D0
    IF (BB.EQ.0.0D0) B(3)=0.0D0
C FOR THE SPECIAL CASE OF THE ENTIRE TRANSFER FUNCTION ON
C ONE TRANSDUCER ONLY, LET hi BE AN IMPULSE (=1).
    IF (K.EQ.0) THEN
            MONE=1
            HONE(1) = 1.0D0
    END IF
C
    IF (GCDUNT.GT.K) GO TO 232
C INCORPORATE THE LINEAR OR QUADRATIC FACTOR INTO THE
C hl POLYNOMIAL
    IF (ZOOUNT. EO. G(GOOUNT)) THEN
            CALL
POLYMULT (MONE, AONE, B, CONE, ZCOUNT, GCOUNT, BB, HONE)
            GCOUNT=GCOUNT+1
            GO TO 5000
    END IF
C INCORPORATE THE LINEAR OR QUADRATIC FACTOR INTO THE
C h2 POLYNOMIAL
232 IF (ZCOUNT.NE.G(GCOUNT)) THEN
            CALL
POLYMULT (MIWO, ATWO, B, CIWO, ZCOUNT, GCOUNT, BB, HIWO)
    END IF
C
5000
            CONTINUE
C SCALE THE COEFFICIENTS TO MAXIMUM OF 1
    HMAX=0.0D0
    DO 96 L=1, MONE
    IF (DABS(HONE(L)).GT.HMAX) HMAX=DABS(HONE(L))
   CONTINUE
    DO 97 L=1.MONE
    HONE (L) =BONE (L) /HPAX
97 CONTINUE
    HMAX=0.0D0
    DO 98 L=1,MIWO
    IF (DABS(HIWO(L)).GT.HMAX) HMAX=DABS(HIWO(L))
98 CONTINUE
    DO 99 L=1,MIWO
```

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```
HIMO(L) =HIMO(L) /HMAX
   CONTINUE
C DETERMINE THE FIGURE OF MERIT FOR THE DESIGN
    IF (K.EO.0) THEN
            HONE(1) = 1.0D0
            FOMOLD=0.0E0
    END IF
    DO 348 L=1,MONE
    HBOTH (L) =HONE (L)
348 CONTINUE
    DO 349 I=1+MONE, MONE+MIWO
    HBOTH (L) =HTWO (L-MONE)
349 CONTINUE
    CALL HSTATS (BOTH, MONE+MIWO, FOM)
C IF THIS IS THE BEST DESIGN RELATIVE TO ALL PAST ONES,
C SAVE THE RESULTS.
    IF (FOM. GT. FOMOLD) THEN
            PRINT *, '** BEST YET FOLLOWS **'
            FOMOLD=FOM
            MONEBEST-MONE
            MIWOBEST=MIWO
            PRINT *, 'FOM=', FOM, ' K=', K
            DO 111 II=1, MONE
            HBESTONE(II) =BONE(II)
111
            CONTINUE
            DO 2222 II=1,MIWO
            HBESTIWO(II) =HIWO(II)
2222
                     CONTINUE
C OPTIONAL CODE TO PROVIDE SPECIFIC INFORMATION CONCERNING
C THE NATURE OF EACH SELECTED ROOT (I.E., PASSBAND OR
C STOPBAND, REAL OR COMPLEX)
            ZCOUNT-1
C
            GCOUNT=1
C
            DO 1234 II=1, DEGREE
            AA=ZEROR(II)
C
            BB=ZEROI(II)
C
            IF (DABS(BB).LT.1.0D-10) BB=0.0D0
            IF (BB.LT.0.0D0) GO TO 1234
C
            IF (GCOUNT.GT.K) GO TO 217
            IF (ZCOUNT. EQ. G(GCOUNT)) THEN
C
                     PRINT *, 'H1 ZERO: ', AA, ' +/-', BB
C
                     AA=DABS(1.0D0-DSQRT(AA*AA+BB*BB))
C
                     IF (AA.LT.1.0D-4) PRINT *, 'STOPBAND'
                     IF (AA.GE.1.0D-4) PRINT *, 'PASSBAND'
C
                     IF (BB.EQ.0.0D0) PRINT *, 'REAL'
                    PRINT *,' '
C
                     ZCOUNT=ZCOUNT+1
```

```
GCOUNT-GCOUNT+1
C
                     GO TO 1234
C
            END IF
C
c 217
                     IF (ZCOUNT. NE. G(GCOUNT)) THEN
                     PRINT *, 'H2 ZERO: ', AA, ' +/-', BB
C
C
                     AA=DABS(1.0D0-DSQRT(AA*AA+BB*BB))
                     IF (AA.LT.1.0D-4) PRINT *, 'STOPBAND'
C
                     IF (AA.GE.1.0D-4) PRINT *, 'PASSBAND'
C
                     IF (BB.EQ.0.0D0) PRINT *, 'REAL'
C
                     PRINT *,' '
C
C
                     ZCOUNT=ZCOUNT+1
            END IF
C
c 1234
                     CONTINUE
    END IF
C
    RETURN
C
    END
C
    SUBROUTINE HSTATS (X, M, FOM)
C THIS SUBROUTINE DETERMINES THE STATISTICS OF THE SAMPLES
C AND RETURNS THE FIGURE OF MERIT (FOM)
    DOUBLE PRECISION X(1024), XMAX, XMIN, XAVG, XVAR
    DOUBLE PRECISION FOM, XSUM, XDUM, XRANGE
    INTEGER M
    XSUM=0.0D0
    XMAX=0.0D0
    XMIN=1.0D20
    XDUM=0.0D0
    DO 10 I=1,M
    XDUM=XDUM+X(I)*X(I)
    XSUM=XSUM+DABS(X(I))
    IF (DABS(X(I)).GT.XMAX) XMAX=DABS(X(I))
    IF (X(I).EQ.0.0D0) GO TO 10
    IF (DABS(X(I)).LT.XMIN) XMIN⇒DABS(X(I))
10 CONTINUE
    XAVG=XSUM/DBLE(M)
    XVAR=(DBLE(M) *XDUM-XSUM*XSUM)/(DBLE(M) * (DBLE(M) -
1.0D0))
    XRANGE=XMAX-XMIN
    IF (XVAR.EQ.0.0D0) XVAR=1.0D-10
    FOM-XAVG/(XRANGE*XVAR)
35 RETURN
    END
C POLYNOMIAL RECONSTITUTION SUBROUTINE. THIS ROUTINE TAKES
C THE LINEAR OR QUADRATIC FACTOR AND MULTIPLIES IT BY THE
CURRENT
C POLINOMIAL.
    SUBROUTINE POLYMULT (M, A, B, C, ZCOUNT, GCOUNT, BB, H)
```

```
DOUBLE PRECISION A(1024), B(3), H(1024)
    DOUBLE PRECISION C(1024), BB
    INTEGER ZCOUNT, M
C IF NO CURRENT POLYNOMIAL YET EXISTS, THEN THE FACTOR
C THE CURRENT FOLYNOMIAL.
    IF (M.EQ.0) THEN
            DO 1002 L=1,3
             A(L) = B(L)
             C(L) = B(L)
1002
                     CONTINUE
             M=1
            GO TO 1421
    END IF
C IF THE CURRENT POLYNOMIAL IS OF ORDER GREATER THAN
THREE, THE
C RECURSIVE RELATIONSHIP APPLIES.
    IF (M.GT.3) THEN
            DO 1001 L=3,M
            C(L) = B(3) *A(L-2)
            C(L) = C(L) + B(2) *A(L-1)
            C(L) = C(L) + B(1) *A(L)
1001
                     CONTINUE
    END IF
C DUE TO VARIABLE ADDRESSING LIMITATIONS, TAKE CARE OF THE
TWO
C HIGH ORDER AND THREE LOWEST ORDER COEFFICIENTS MANUALLY
    C(M+2) = B(3) *A(M)
    C(M+1)=B(3)*A(M-1)+B(2)*A(M)
    C(3) = B(3) *A(1) +B(2) *A(2) +B(1) *A(3)
    C(2)=B(2)*A(1)+B(1)*A(2)
    C(1) = B(1) *A(1)
C UPDATE THE NUMBER OF ZEROS USED.
1421
            ZCOUNT=ZCOUNT+1
C ORDER INCREASES BY TWO FOR A QUADRATIC FACTOR, ONE FOR A
C LINEAR FACTOR.
    IF (BB. NE. 0.0D0) M=M+2
    IF (BB. EQ. 0.0D0) M=M+1
C ESTABLISH THE NEW CURRENT POLYNOMIAL
C AND THE NEW IMPULSE RESPONSE
    DO 131 L=1,M
```

A(L) =C(L) H(L) =C(L) 131 CONTINUE RETURN END

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